

Instrument ID: **MSVOA_L**

Daily Analysis Runlog For Sequence/QC Batch ID # VL121824

Review By	Semsettin Yesilyurt	Review On	12/19/2024 1:32:59 PM		
Supervise By	Maresh Dadoda	Supervise On	12/19/2024 2:45:49 PM		
SubDirectory	VL121824	HP Acquire Method	TO15AIR.M	HP Processing Method	VL121624AIR.M
STD. NAME		STD REF.#			
Tune/Reschk		AP2556			
Initial Calibration Stds		AP2560,AP2561,AP2562			
CCC		AP2560,MDL,AP2561,AP2562			
Internal Standard/PEM		AP2556			
ICV/I.BLK		AP2558			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VL041787.D	18 Dec 2024 08:06	SY/MD	Ok
2	VSTDCCC010	VL041788.D	18 Dec 2024 10:29	SY/MD	Ok,M
3	VL1218ABL01	VL041789.D	18 Dec 2024 11:57	SY/MD	Ok
4	VL1218ABL02	VL041790.D	18 Dec 2024 12:28	SY/MD	Ok
5	MDL 0.4	VL041791.D	18 Dec 2024 12:59	SY/MD	Ok,M
6	MDL 0.4	VL041792.D	18 Dec 2024 13:30	SY/MD	Ok,M
7	MDL 0.03	VL041793.D	18 Dec 2024 14:32	SY/MD	Ok,M
8	MDL 0.03	VL041794.D	18 Dec 2024 15:03	SY/MD	Ok,M
9	VL1218ABS01	VL041795.D	18 Dec 2024 16:06	SY/MD	Ok,M
10	QC121724 CAN 10158	VL041796.D	18 Dec 2024 16:37	SY/MD	Ok
11	P5293-01	VL041797.D	18 Dec 2024 17:10	SY/MD	Ok,M
12	P5293-01DUP	VL041798.D	18 Dec 2024 17:43	SY/MD	Ok,M
13	P5293-02	VL041799.D	18 Dec 2024 18:17	SY/MD	Ok,M
14	P5293-03	VL041800.D	18 Dec 2024 18:50	SY/MD	Dilution
15	P5296-01	VL041801.D	18 Dec 2024 19:24	SY/MD	Ok,M
16	P5294-01	VL041802.D	18 Dec 2024 19:58	SY/MD	Ok,M
17	P5294-02	VL041803.D	18 Dec 2024 20:32	SY/MD	Ok,M
18	P5294-03	VL041804.D	18 Dec 2024 21:03	SY/MD	Not Ok
19	P5292-01	VL041805.D	18 Dec 2024 21:35	SY/MD	Not Ok
20	P5292-02	VL041806.D	18 Dec 2024 22:06	SY/MD	Not Ok
21	P5295-01	VL041807.D	18 Dec 2024 22:37	SY/MD	Dilution

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CCC	AP2560,MDL,AP2561,AP2562				
Internal Standard/PEM	AP2556				
ICV/I.BLK	AP2558				
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

22	P5295-02DL	VL041808.D	18 Dec 2024 23:08	SY/MD	Ok,M
23	VIBLK	VL041809.D	18 Dec 2024 23:40	SY/MD	Ok
24	P5293-03DL	VL041810.D	19 Dec 2024 00:11	SY/MD	Ok,M
25	P5294-03	VL041811.D	19 Dec 2024 00:44	SY/MD	Ok,M
26	P5292-01	VL041812.D	19 Dec 2024 01:18	SY/MD	Ok,M
27	P5292-02	VL041813.D	19 Dec 2024 01:55	SY/MD	Ok,M
28	P5295-01DL	VL041814.D	19 Dec 2024 02:26	SY/MD	Ok,M
29	P5295-02	VL041815.D	19 Dec 2024 02:59	SY/MD	Dilution
30	VSTDCCC010	VL041816.D	19 Dec 2024 03:31	SY/MD	Ok

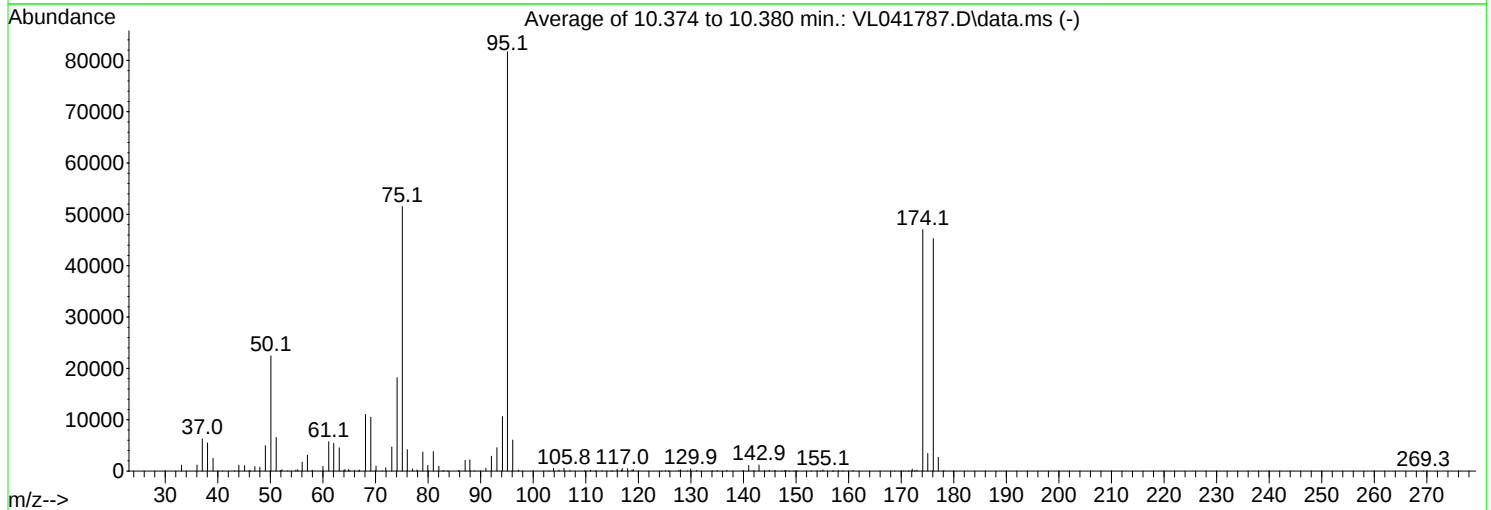
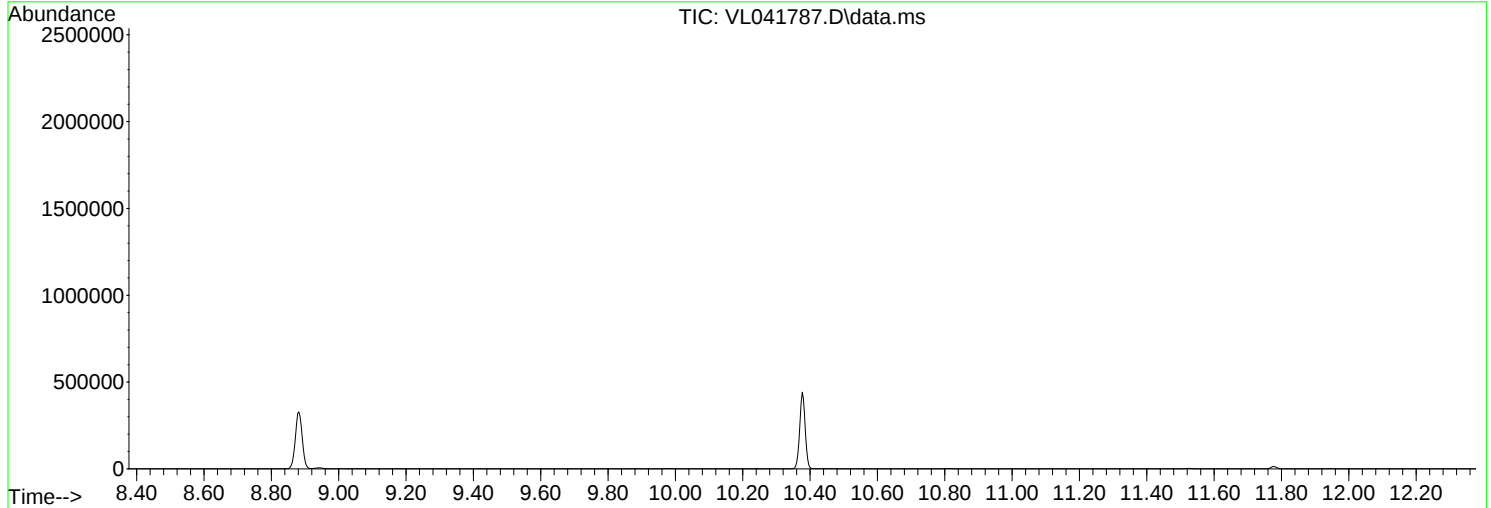
M : Manual Integration

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
Data File : VL041787.D
Acq On : 18 Dec 2024 08:06
Operator : SY/MD
Sample : BFB
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
Last Update : Thu Dec 19 07:32:09 2024



AutoFind: Scans 3178, 3179, 3180; Background Corrected with Scan 3165

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	27.5	22445	PASS
75	95	30	66	63.1	51533	PASS
95	95	100	100	100.0	81685	PASS
96	95	5	9	7.4	6040	PASS
173	174	0.00	2	0.4	189	PASS
174	95	50	120	57.6	47024	PASS
175	174	4	9	7.3	3435	PASS
176	174	93	101	96.3	45299	PASS
177	176	5	9	5.9	2683	PASS

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041788.D
 Acq On : 18 Dec 2024 10:29
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDCCC010

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 12/19/2024
 Supervised By : Mahesh Dadoda 12/19/2024

Quant Time: Dec 19 07:49:00 2024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Thu Dec 19 07:32:09 2024

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.784	49	76040	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.952	114	198370	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.881	117	172190	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.377 95 150705 10.503 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 105.000%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.495	85	112841	8.340	ppbv	96
3) Chlorodifluoromethane	1.473	51	112096	9.040	ppbv	98
4) Chloromethane	1.531	50	48418	9.146	ppbv	99
5) Vinyl Chloride	1.580	62	46631	8.743	ppbv	99
6) Bromomethane	1.667	94	22202	8.821	ppbv	98
7) Chloroethane	1.703	64	19203	9.269	ppbv	99
8) Dichlorotetrafluoroethane	1.554	85	111653	9.220	ppbv	96
9) Propene	1.483	41	41427	7.708	ppbv	96
10) Heptane	4.975	43	111650	8.087	ppbv	96
11) Trichlorofluoromethane	1.874	101	134331	9.347	ppbv	99
12) 1,1,2-Trichlorotrifluo...	2.136	101	96469	8.825	ppbv	98
13) Ethanol	1.719	45	3276m	7.361	ppbv	
14) Bromoethene	1.780	108	33118	8.974	ppbv	97
15) Acetone	1.832	43	111294	7.522	ppbv	99
16) 1,3-Butadiene	1.606	39	53299	9.149	ppbv	97
17) tert-Butyl alcohol	2.036	59	135516	9.233	ppbv	100
18) 1,1-Dichloroethene	2.033	96	41224	8.996	ppbv	93
19) Isopropyl Alcohol	1.884	45	64831	7.303	ppbv	99
20) Methylene Chloride	2.059	84	35699	8.462	ppbv	96
21) Allyl Chloride	2.094	41	74468	9.281	ppbv	99
22) trans-1,2-Dichloroethene	2.337	96	45331	8.641	ppbv	94
23) Vinyl Acetate	2.460	43	52709	7.673	ppbv #	95
24) 1,1-Dichloroethane	2.405	63	95779	9.020	ppbv	99
25) Ethyl Acetate	2.832	43	238771	8.675	ppbv	100
26) Hexane	2.822	57	87593	8.386	ppbv	95
27) Carbon Disulfide	2.149	76	85318	9.064	ppbv	98
28) Methyl tert-Butyl Ether	2.437	73	67965	8.021	ppbv	97
29) Chloroform	2.845	83	145536	9.143	ppbv	97
30) Cyclohexane	3.855	84	74346	8.366	ppbv	97
31) cis-1,2-Dichloroethene	2.719	61	92460	8.567	ppbv	98
32) 1,1,1-Trichloroethane	3.366	97	144038	8.902	ppbv	99
34) 2-Butanone	2.554	43	139509	8.742	ppbv	99
35) Carbon Tetrachloride	3.761	117	147726	10.490	ppbv	99
36) Benzene	3.654	78	178114	8.680	ppbv	98
37) 1,2-Dichloroethane	3.214	62	121412	9.509	ppbv	100
38) Trichloroethene	4.548	130	70175	8.057	ppbv	98
39) 1,2-Dichloropropane	4.308	63	64154	8.645	ppbv	97
40) 1,4-Dioxane	4.583	88	30262	8.347	ppbv #	96
41) Tetrahydrofuran	3.052	42	74734	8.359	ppbv	98
42) Bromodichloromethane	4.486	83	140832	9.588	ppbv	95
43) Methyl Methacrylate	4.878	69	69545	8.895	ppbv	99
44) 2,2,4-Trimethylpentane	4.635	57	302332	8.627	ppbv	97
45) t-1,3-Dichloropropene	6.412	75	69601	8.607	ppbv	96

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041788.D
 Acq On : 18 Dec 2024 10:29
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDCCC010

Quant Time: Dec 19 07:49:00 2024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Thu Dec 19 07:32:09 2024

Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 12/19/2024
 Supervised By : Mahesh Dadoda 12/19/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.593	75	89328m	8.701	ppbv	
47) 1,1,2-Trichloroethane	6.580	97	67717	8.736	ppbv	93
48) Dibromochloromethane	7.386	129	101875	9.625	ppbv	97
49) Bromoform	9.483	173	83649	9.655	ppbv	99
50) 4-Methyl-2-Pentanone	5.752	43	186865	8.600	ppbv	99
51) 2-Hexanone	7.438	43	149217	8.472	ppbv #	98
52) Tetrachloroethene	8.224	164	59756	7.824	ppbv	93
53) Toluene	6.930	91	195158	8.389	ppbv	99
54) 1,2-Dibromoethane	7.648	107	94138	8.814	ppbv	97
56) 1,1,1,2-Tetrachloroethane	8.924	131	87869	9.626	ppbv	99
57) Chlorobenzene	8.920	112	157039	8.912	ppbv	97
58) Ethyl Benzene	9.354	91	281290	9.074	ppbv	97
59) m/p-Xylene	9.548	91	466737m	18.725	ppbv	
60) o-Xylene	9.963	91	232028	9.283	ppbv	100
61) Styrene	9.869	104	103216	9.550	ppbv	99
62) Isopropylbenzene	10.539	105	352903	9.397	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.963	83	144197	9.216	ppbv	99
64) n-propylbenzene	10.989	120	90043	9.483	ppbv	99
65) tert-Butylbenzene	11.532	119	314940	9.295	ppbv	98
66) Benzyl Chloride	11.616	91	22164	7.426	ppbv	100
67) sec-Butylbenzene	11.769	105	442201	9.431	ppbv	100
69) p-Isopropyltoluene	11.927	119	369746	9.569	ppbv	99
70) n-Butylbenzene	12.283	91	373294	9.999	ppbv	99
71) 2-Chlorotoluene	10.901	91	264866	9.240	ppbv	100
72) 4-Ethyltoluene	11.124	105	285469	9.676	ppbv	99
73) 1,3,5-Trimethylbenzene	11.202	105	242783	9.918	ppbv	97
74) 1,2,4-Trimethylbenzene	11.536	105	263301	10.003	ppbv	90
75) 1,3-Dichlorobenzene	11.607	146	154857	9.227	ppbv	99
76) 1,4-Dichlorobenzene	11.671	146	154115	9.296	ppbv	99
77) 1,2-Dichlorobenzene	11.947	146	145375	9.002	ppbv	99
78) Hexachloro-1,3-Butadiene	13.882	225	93437	7.804	ppbv	100
79) Naphthalene	13.513	128	200311	10.921	ppbv	98
80) Naphthalene,2-methyl-	14.458	142	94903	16.728	ppbv	97
81) 1,2,4-Trichlorobenzene	13.442	180	93675	9.346	ppbv	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

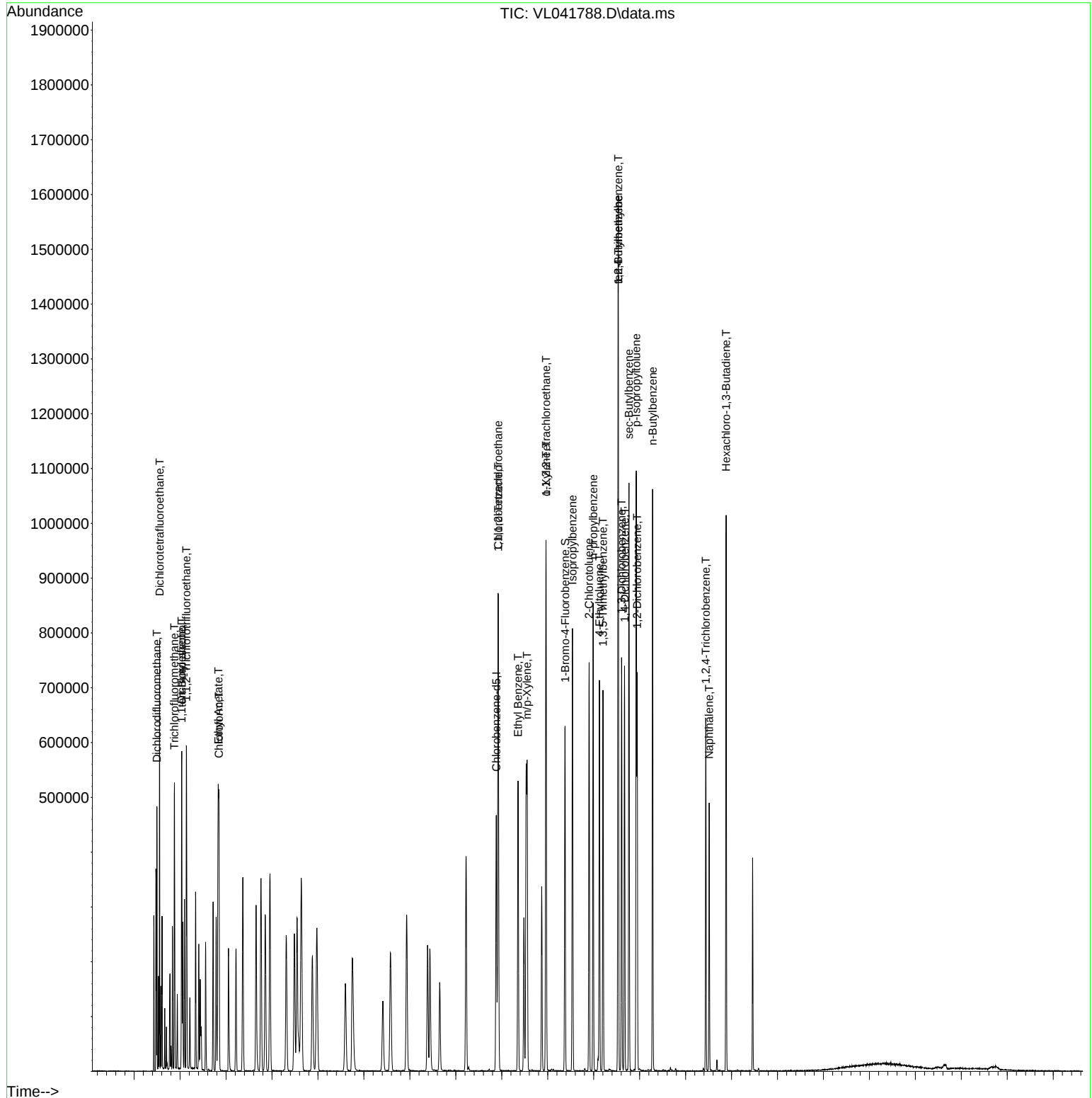
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
Data File : VL041788.D
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Operator : SY/MD
Sample : VSTDCCC010
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VSTDCCC010

Manual Integrations
APPROVED

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Supervised By : Mahesh Dadoda 12/19/2024

Quant Time: Dec 19 07:49:00 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Thu Dec 19 07:32:09 2024
Response via : Initial Calibration



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 QLast Update : Thu Dec 19 07:32:09 2024
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	111	0.00
2 T	Dichlorodifluoromethane	1.779	1.484	16.6	98	0.00
3	Chlorodifluoromethane	1.631	1.474	9.6	104	0.00
4	Chloromethane	0.696	0.637	8.5	112	0.00
5 T	Vinyl Chloride	0.701	0.613	12.6	109	0.00
6 T	Bromomethane	0.331	0.292	11.8	110	0.00
7	Chloroethane	0.272	0.253	7.0	110	0.00
8 T	Dichlorotetrafluoroethane	1.593	1.468	7.8	111	0.00
9 T	Propene	0.707	0.545	22.9	95	0.00
10 T	Heptane	1.816	1.468	19.2	95	0.00
11 T	Trichlorofluoromethane	1.890	1.767	6.5	109	0.00
12 T	1,1,2-Trichlorotrifluoroeth	1.437	1.269	11.7	105	0.00
13	Ethanol	0.059	0.043#	27.1	132	0.00
14 T	Bromoethene	0.485	0.436	10.1	102	0.00
15 T	Acetone	1.946	1.464	24.8	111	0.00
16 T	1,3-Butadiene	0.766	0.701	8.5	113	0.00
17	tert-Butyl alcohol	1.930	1.782	7.7	108	0.00
18 T	1,1-Dichloroethene	0.603	0.542	10.1	106	0.00
19 T	Isopropyl Alcohol	1.167	0.853	26.9	110	0.00
20 T	Methylene Chloride	0.555	0.469	15.5	109	0.00
21 T	Allyl Chloride	1.055	0.979	7.2	107	0.00
22 T	trans-1,2-Dichloroethene	0.690	0.596	13.6	103	0.00
23 T	Vinyl Acetate	0.903	0.693	23.3	97	0.00
24 T	1,1-Dichloroethane	1.396	1.260	9.7	107	0.00
25 T	Ethyl Acetate	3.620	3.140	13.3	101	0.00
26 T	Hexane	1.374	1.152	16.2	98	0.00
27 T	Carbon Disulfide	1.238	1.122	9.4	92	0.00
28 T	Methyl tert-Butyl Ether	1.114	0.894	19.7	90	0.00
29 T	Chloroform	2.093	1.914	8.6	108	0.00
30 T	Cyclohexane	1.169	0.978	16.3	99	0.00
31 T	cis-1,2-Dichloroethene	1.419	1.216	14.3	101	0.00
32 T	1,1,1-Trichloroethane	2.128	1.894	11.0	104	0.00
33 I	1,4-Difluorobenzene	1.000	1.000	0.0	108	0.00
34 T	2-Butanone	0.805	0.703	12.7	103	0.00
35 T	Carbon Tetrachloride	0.710	0.745	-4.9	106	0.00
36 T	Benzene	1.034	0.898	13.2	99	0.00
37 T	1,2-Dichloroethane	0.644	0.612	5.0	109	0.00
38 T	Trichloroethene	0.439	0.354	19.4	98	0.00
39 T	1,2-Dichloropropane	0.374	0.323	13.6	99	0.00
40 T	1,4-Dioxane	0.183	0.153	16.4	99	0.00
41 T	Tetrahydrofuran	0.451	0.377	16.4	95	0.00
42 T	Bromodichloromethane	0.740	0.710	4.1	99	0.00
43	Methyl Methacrylate	0.394	0.351	10.9	96	0.00
44 T	2,2,4-Trimethylpentane	1.767	1.524	13.8	98	0.00
45 T	t-1,3-Dichloropropene	0.408	0.351	14.0	90	0.00
46 T	cis-1,3-Dichloropropene	0.518	0.450	13.1	91	0.00
47 T	1,1,2-Trichloroethane	0.391	0.341	12.8	100	0.00
48 T	Dibromochloromethane	0.534	0.514	3.7	94	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
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 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 LabSampleId :
 VSTDCCC010

Quant Time: Dec 19 07:49:00 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Thu Dec 19 07:32:09 2024
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	Bromoform	0.437	0.422	3.4	91	0.00
50 T	4-Methyl-2-Pentanone	1.095	0.942	14.0	96	0.00
51 T	2-Hexanone	0.888	0.752	15.3	95	0.00
52 T	Tetrachloroethene	0.385	0.301	21.8	95	0.00
53 T	Toluene	1.173	0.984	16.1	95	0.00
54 T	1,2-Dibromoethane	0.538	0.475	11.7	100	0.00
55 I	Chlorobenzene-d5	1.000	1.000	0.0	105	0.00
56	1,1,1,2-Tetrachloroethane	0.530	0.510	3.8	101	0.00
57 T	Chlorobenzene	1.023	0.912	10.9	101	0.00
58 T	Ethyl Benzene	1.800	1.634	9.2	97	0.00
59 T	m/p-Xylene	1.448	1.355	6.4	101	0.00
60 T	o-Xylene	1.452	1.348	7.2	100	0.00
61 T	Styrene	0.628	0.599	4.6	94	0.00
62	Isopropylbenzene	2.181	2.049	6.1	101	0.00
63 T	1,1,2,2-Tetrachloroethane	0.909	0.837	7.9	105	0.00
64	n-propylbenzene	0.551	0.523	5.1	103	0.00
65	tert-Butylbenzene	1.968	1.829	7.1	102	0.00
66 T	Benzyl Chloride	0.173	0.129	25.4	71	0.00
67	sec-Butylbenzene	2.723	2.568	5.7	103	0.00
68 S	1-Bromo-4-Fluorobenzene	0.833	0.875	-5.0	109	0.00
69	p-Isopropyltoluene	2.244	2.147	4.3	102	0.00
70	n-Butylbenzene	2.168	2.168	0.0	104	0.00
71	2-Chlorotoluene	1.665	1.538	7.6	100	0.00
72 T	4-Ethyltoluene	1.713	1.658	3.2	101	0.00
73 T	1,3,5-Trimethylbenzene	1.422	1.410	0.8	101	0.00
74 T	1,2,4-Trimethylbenzene	1.529	1.529	0.0	103	0.00
75 T	1,3-Dichlorobenzene	0.975	0.899	7.8	102	0.00
76 T	1,4-Dichlorobenzene	0.963	0.895	7.1	101	0.00
77 T	1,2-Dichlorobenzene	0.938	0.844	10.0	103	0.00
78 T	Hexachloro-1,3-Butadiene	0.695	0.543	21.9	98	0.00
79 T	Naphthalene	1.065	1.163	-9.2	97	0.00
80 T	Naphthalene,2-methyl-	0.329	0.551	-67.5#	89	0.00
81 T	1,2,4-Trichlorobenzene	0.582	0.544	6.5	92	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041790.D
 Acq On : 18 Dec 2024 12:28
 Operator : SY/MD
 Sample : VL1218ABL02
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL1218ABL02

Quant Time: Dec 19 07:52:46 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Thu Dec 19 07:32:09 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.793	49	70249	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.965	114	175356	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.891	117	151060	10.000	ppbv	0.01

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	10.383	95	127683	10.143	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	101.400%

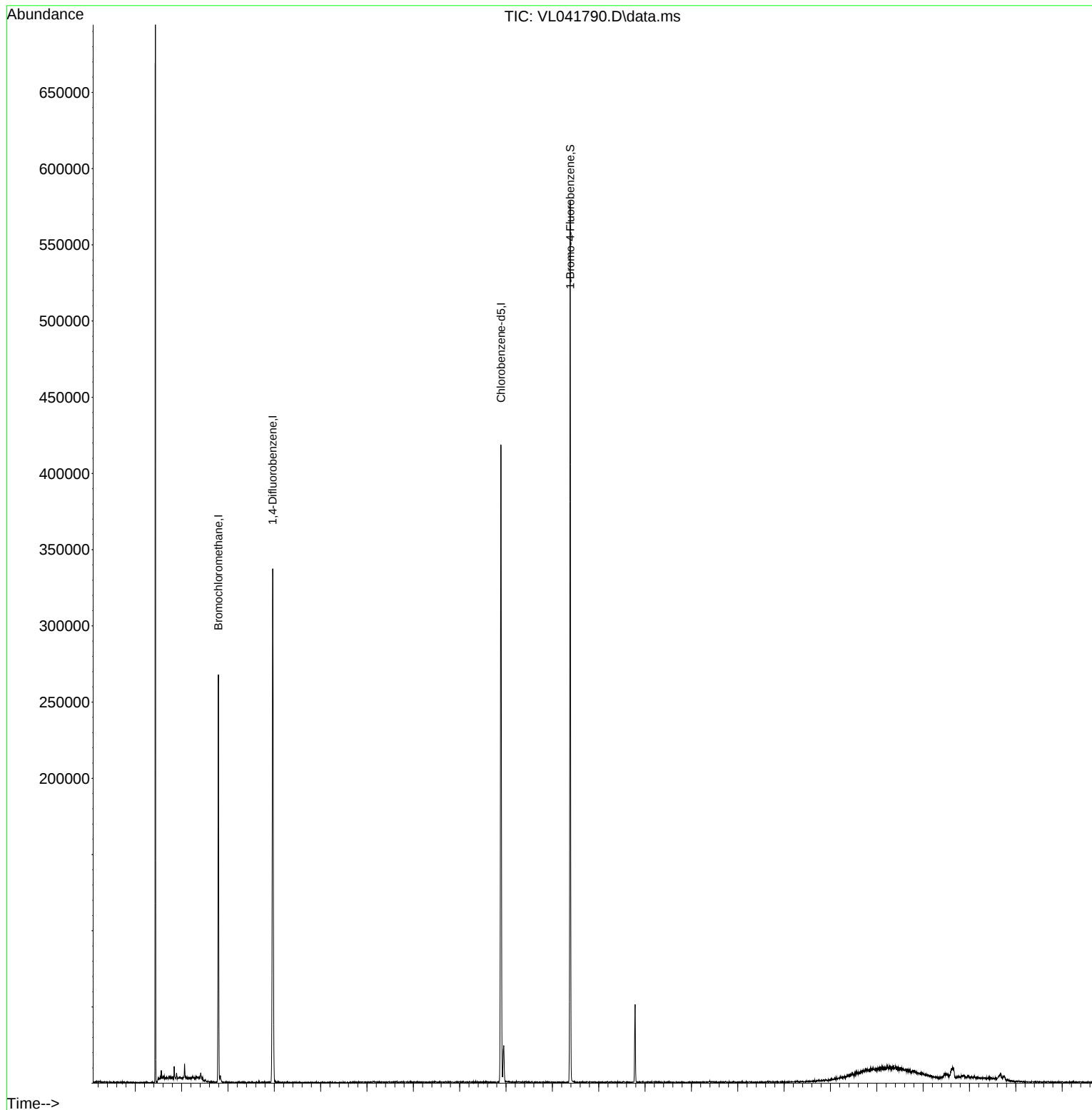
Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
Data File : VL041790.D
Acq On : 18 Dec 2024 12:28
Operator : SY/MD
Sample : VL1218ABL02
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VL1218ABL02

Quant Time: Dec 19 07:52:46 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
QLast Update : Thu Dec 19 07:32:09 2024
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041795.D
 Acq On : 18 Dec 2024 16:06
 Operator : SY/MD
 Sample : VL1218ABS01
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL1218ABS01

Quant Time: Dec 19 07:55:13 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2
 QLast Update : Thu Dec 19 07:32:09 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 12/19/2024
 Supervised By : Mahesh Dadoda 12/19/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.797	49	72600	10.000	ppbv	0.01
33) 1,4-Difluorobenzene	3.972	114	183255	10.000	ppbv	0.02
55) Chlorobenzene-d5	8.898	117	159496	10.000	ppbv	0.02

System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.390	95	143879	10.825	ppbv	0.02
Spiked Amount	10.000	Range	65 - 135	Recovery	=	108.300%

Target Compounds	Qvalue					
2) Dichlorodifluoromethane	1.502	85	112657	8.721	ppbv	99
3) Chlorodifluoromethane	1.479	51	102210	8.633	ppbv	97
4) Chloromethane	1.538	50	46901	9.279	ppbv	100
5) Vinyl Chloride	1.583	62	44104	8.661	ppbv	98
6) Bromomethane	1.674	94	21119	8.788	ppbv	99
7) Chloroethane	1.709	64	18047	9.123	ppbv	98
8) Dichlorotetrafluoroethane	1.557	85	110410	9.549	ppbv	93
9) Propene	1.486	41	35133	6.846	ppbv	98
10) Heptane	5.001	43	106572	8.085	ppbv	100
11) Trichlorofluoromethane	1.884	101	132171	9.632	ppbv	98
12) 1,1,2-Trichlorotrifluo...	2.146	101	96243	9.222	ppbv	95
13) Ethanol	1.725	45	2865	6.742	ppbv	98
14) Bromoethene	1.787	108	30065	8.533	ppbv	96
15) Acetone	1.842	43	110277	7.807	ppbv	97
16) 1,3-Butadiene	1.612	39	52562	9.450	ppbv	92
17) tert-Butyl alcohol	2.046	59	122872	8.768	ppbv	100
18) 1,1-Dichloroethene	2.039	96	38415	8.780	ppbv	80
19) Isopropyl Alcohol	1.890	45	59665	7.039	ppbv #	37
20) Methylene Chloride	2.068	84	35310	8.766	ppbv	94
21) Allyl Chloride	2.104	41	68525	8.945	ppbv	95
22) trans-1,2-Dichloroethene	2.350	96	42821	8.550	ppbv	98
23) Vinyl Acetate	2.470	43	43463	6.627	ppbv	97
24) 1,1-Dichloroethane	2.418	63	92680	9.142	ppbv	99
25) Ethyl Acetate	2.845	43	223419	8.502	ppbv	99
26) Hexane	2.835	57	81413	8.164	ppbv	93
27) Carbon Disulfide	2.159	76	76433	8.505	ppbv	94
28) Methyl tert-Butyl Ether	2.447	73	64037	7.915	ppbv	99
29) Chloroform	2.858	83	139818	9.200	ppbv	99
30) Cyclohexane	3.871	84	66237	7.806	ppbv	98
31) cis-1,2-Dichloroethene	2.732	61	82779	8.033	ppbv	99
32) 1,1,1-Trichloroethane	3.379	97	136055	8.807	ppbv	98
34) 2-Butanone	2.564	43	125687	8.525	ppbv	100
35) Carbon Tetrachloride	3.777	117	141202	10.854	ppbv	97
36) Benzene	3.667	78	164423	8.674	ppbv	98
37) 1,2-Dichloroethane	3.230	62	113600	9.631	ppbv	99
38) Trichloroethene	4.564	130	63280	7.864	ppbv	93
39) 1,2-Dichloropropane	4.331	63	58553	8.541	ppbv	92
40) 1,4-Dioxane	4.599	88	28884	8.624	ppbv #	92
41) Tetrahydrofuran	3.065	42	67017	8.114	ppbv	99
42) Bromodichloromethane	4.509	83	133513	9.839	ppbv	95
43) Methyl Methacrylate	4.897	69	61014	8.448	ppbv	99
44) 2,2,4-Trimethylpentane	4.658	57	284140	8.777	ppbv	96
45) t-1,3-Dichloropropene	6.438	75	58865	7.879	ppbv	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041795.D
 Acq On : 18 Dec 2024 16:06
 Operator : SY/MD
 Sample : VL1218ABS01
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL1218ABS01

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 12/19/2024
 Supervised By : Mahesh Dadoda 12/19/2024

Quant Time: Dec 19 07:55:13 2024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Thu Dec 19 07:32:09 2024

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.625	75	74198	7.824	ppbv	89
47) 1,1,2-Trichloroethane	6.603	97	64954	9.071	ppbv	93
48) Dibromochloromethane	7.406	129	96996	9.920	ppbv	99
49) Bromoform	9.500	173	80840	10.100	ppbv	99
50) 4-Methyl-2-Pentanone	5.774	43	175817	8.759	ppbv	99
51) 2-Hexanone	7.457	43	142403	8.752	ppbv #	96
52) Tetrachloroethene	8.244	164	54531	7.729	ppbv	93
53) Toluene	6.952	91	177625	8.265	ppbv	99
54) 1,2-Dibromoethane	7.671	107	90389	9.161	ppbv	92
56) 1,1,1,2-Tetrachloroethane	8.943	131	86262	10.202	ppbv	100
57) Chlorobenzene	8.937	112	153659	9.415	ppbv	99
58) Ethyl Benzene	9.370	91	267795	9.326	ppbv	98
59) m/p-Xylene	9.564	91	455916m	19.747	ppbv	
60) o-Xylene	9.975	91	229919	9.930	ppbv	100
61) Styrene	9.885	104	95536	9.543	ppbv	100
62) Isopropylbenzene	10.552	105	339599	9.762	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.975	83	142272	9.817	ppbv	100
64) n-propylbenzene	11.002	120	86896	9.880	ppbv	97
65) tert-Butylbenzene	11.542	119	307835	9.808	ppbv	97
66) Benzyl Chloride	11.626	91	21939	7.936	ppbv	97
67) sec-Butylbenzene	11.778	105	441578	10.167	ppbv	99
69) p-Isopropyltoluene	11.937	119	366111	10.229	ppbv	99
70) n-Butylbenzene	12.293	91	366451	10.597	ppbv	98
71) 2-Chlorotoluene	10.914	91	255490	9.622	ppbv	99
72) 4-Ethyltoluene	11.137	105	277025	10.137	ppbv	99
73) 1,3,5-Trimethylbenzene	11.215	105	240648	10.613	ppbv	100
74) 1,2,4-Trimethylbenzene	11.548	105	260378	10.679	ppbv	96
75) 1,3-Dichlorobenzene	11.620	146	152493	9.809	ppbv	98
76) 1,4-Dichlorobenzene	11.684	146	152432	9.926	ppbv	97
77) 1,2-Dichlorobenzene	11.960	146	148412	9.922	ppbv	99
78) Hexachloro-1,3-Butadiene	13.892	225	88582	7.987	ppbv	100
79) Naphthalene	13.523	128	189506	11.154	ppbv	99
80) Naphthalene,2-methyl-	14.468	142	82155	15.634	ppbv	96
81) 1,2,4-Trichlorobenzene	13.452	180	89052	9.592	ppbv	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

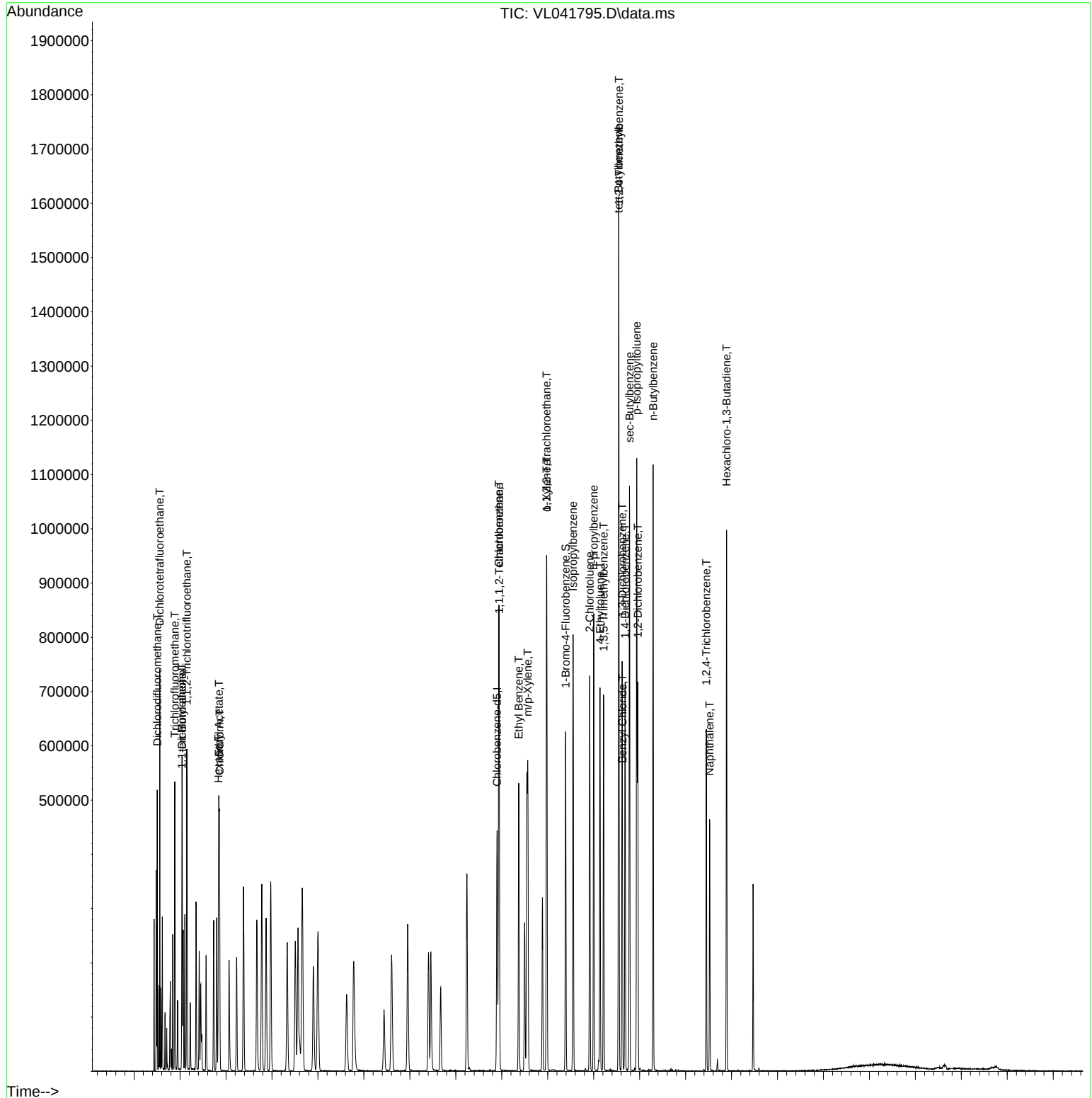
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
Data File : VL041795.D
Acq On : 18 Dec 2024 16:06
Operator : SY/MD
Sample : VL1218ABS01
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VL1218ABS01

Manual Integrations
APPROVED

Reviewed By :Semsettin Yesilyurt 12/19/2024
Supervised By :Mahesh Dadoda 12/19/2024

Quant Time: Dec 19 07:55:13 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LFri Aug
QLast Update : Thu Dec 19 07:32:09 2024
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
 Data File : VL041796.D
 Acq On : 18 Dec 2024 16:37
 Operator : SY/MD
 Sample : QC121724 CAN 10158
 Misc : 400mL/MSVOA_L
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 QC121724 CAN 10158

Quant Time: Dec 19 07:56:01 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Thu Dec 19 07:32:09 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.793	49	73011	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.968	114	179409	10.000	ppbv	0.01
55) Chlorobenzene-d5	8.891	117	152308	10.000	ppbv	0.01

System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.386	95	125911	9.920	ppbv	0.01
Spiked Amount	10.000	Range	65 - 135	Recovery	=	99.200%

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL121824\
Data File : VL041796.D
Acq On : 18 Dec 2024 16:37
Operator : SY/MD
Sample : QC121724 CAN 10158
Misc : 400mL/MSVOA_L
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
QC121724 CAN 10158

Quant Time: Dec 19 07:56:01 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL121624AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
QLast Update : Thu Dec 19 07:32:09 2024
Response via : Initial Calibration

