

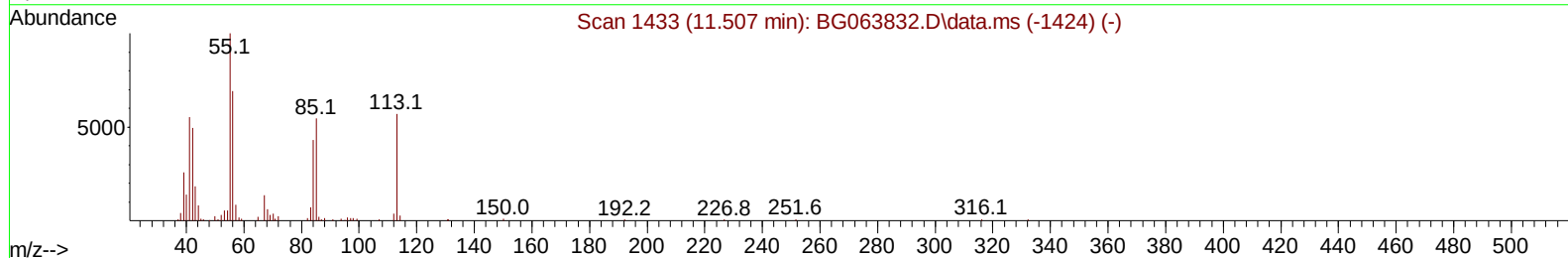
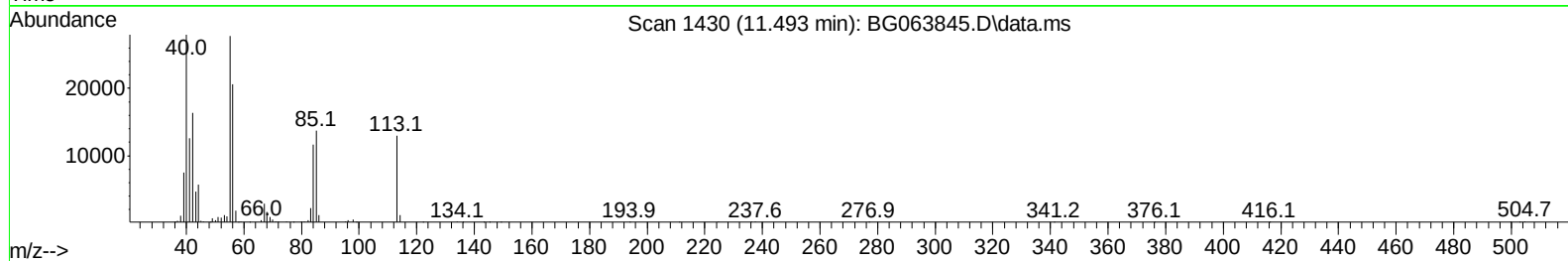
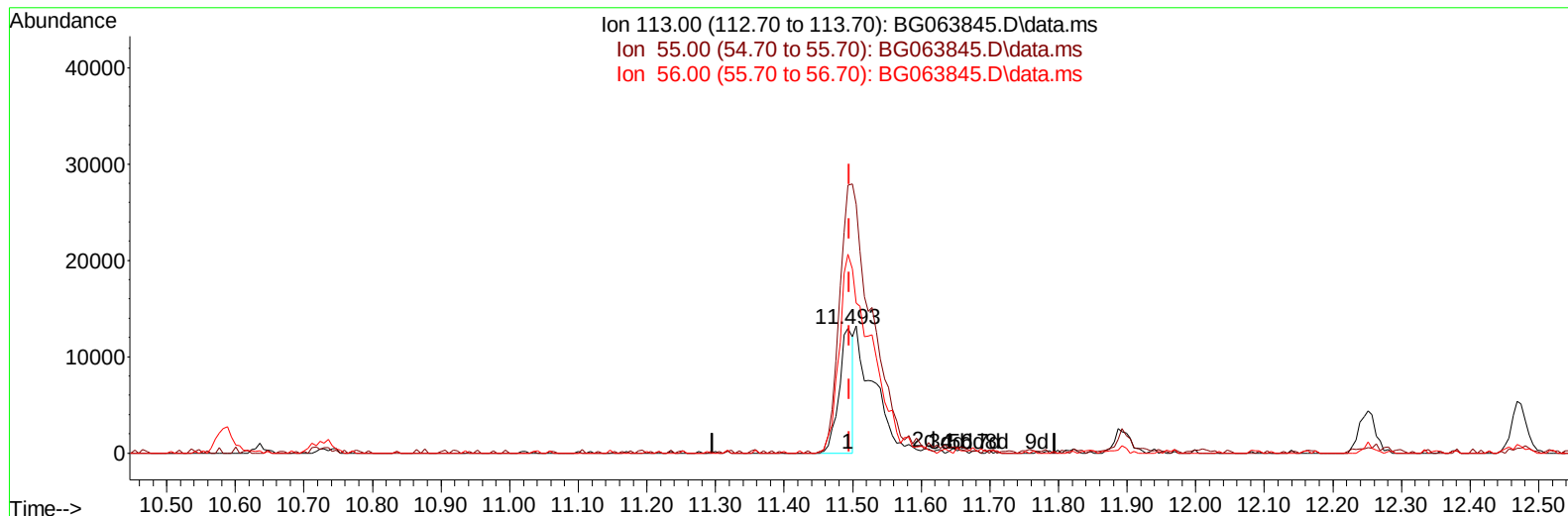
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063845.D
 Acq On : 26 Dec 2024 22:02
 Operator : RC/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 27 01:24:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/27/2024
 Supervised By :mohammad ahmed 12/30/2024



TIC: BG063845.D\data.ms

(34) Caprolactam

11.493min (-0.002) 9.58 ng/u1

response 18300

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	202.90	214.55
56.00	133.90	159.00
0.00	0.00	0.00

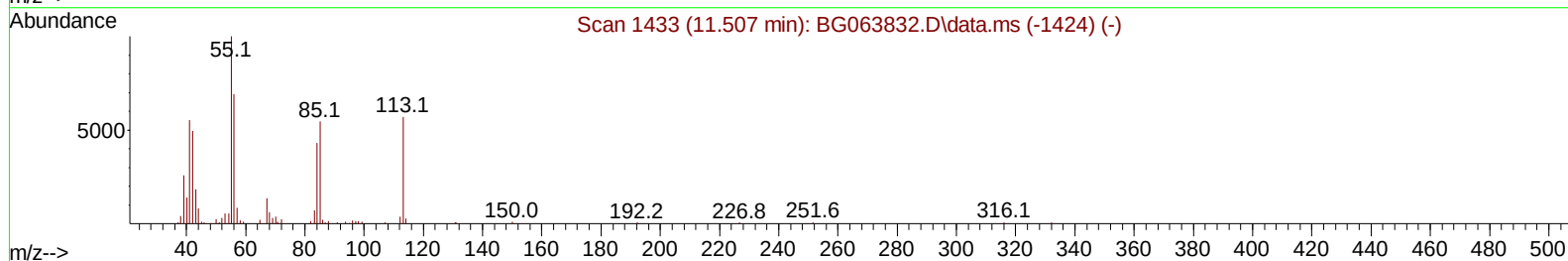
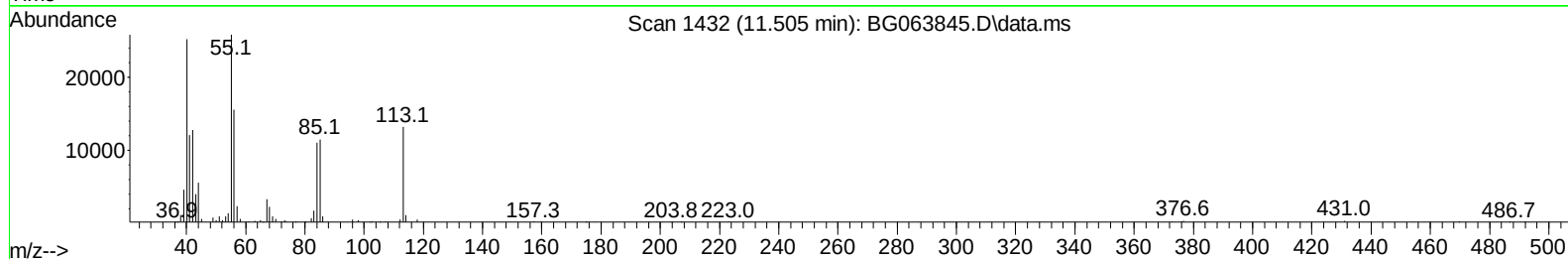
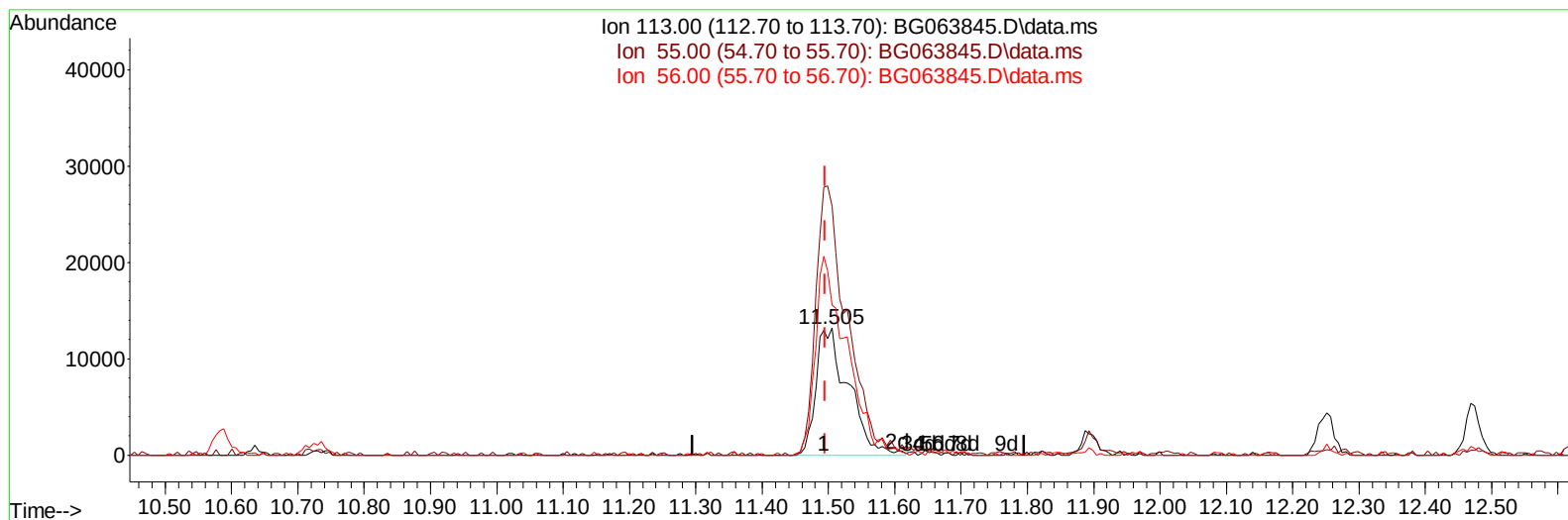
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TIC: BG063845.D\data.ms

(34) Caprolactam

11.505min (+ 0.010) 23.29 ng/ul m

response 44472

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	202.90	195.72
56.00	133.90	117.93
0.00	0.00	0.00

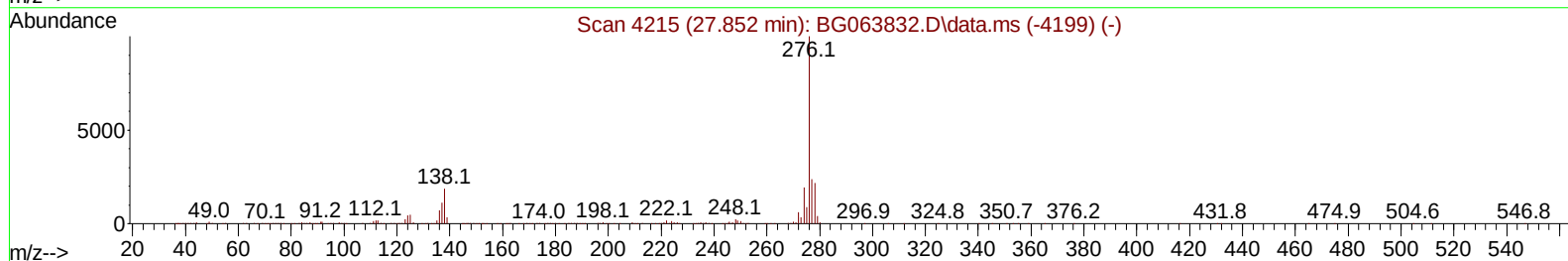
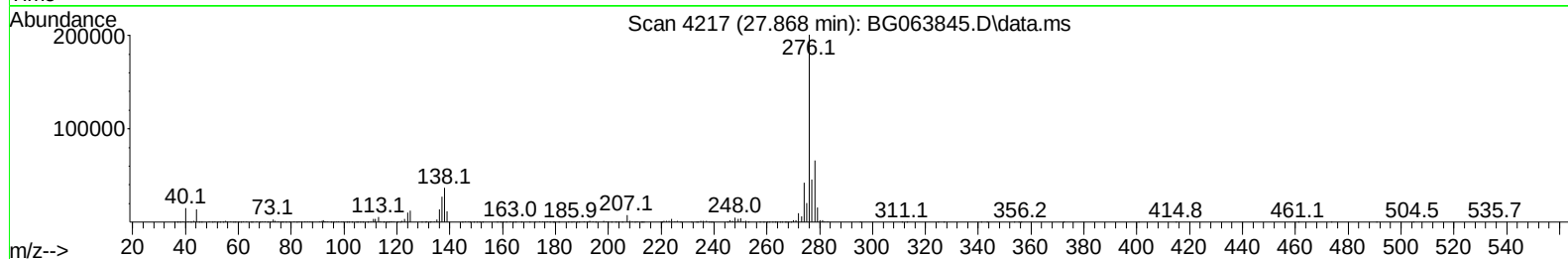
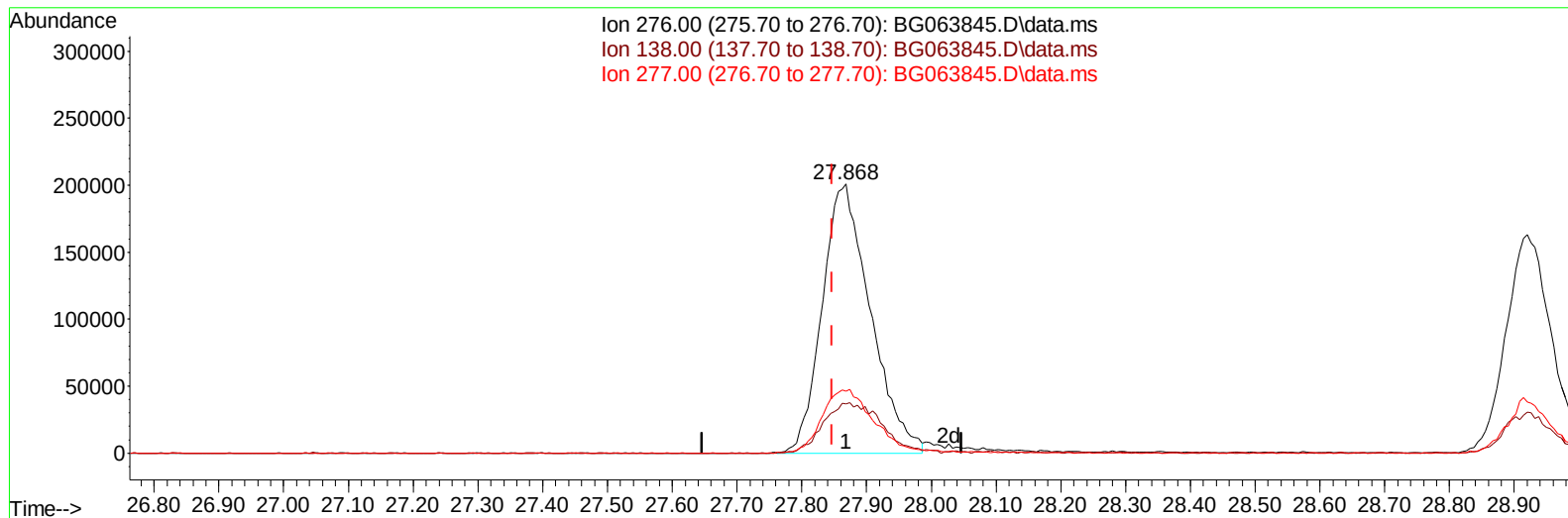
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Manual IntegrationsAPPROVED

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TIC: BG063845.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

27.868min (+ 0.021) 21.31 ng/ul

response 1037816

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	17.10	18.29
277.00	23.00	22.91
0.00	0.00	0.00

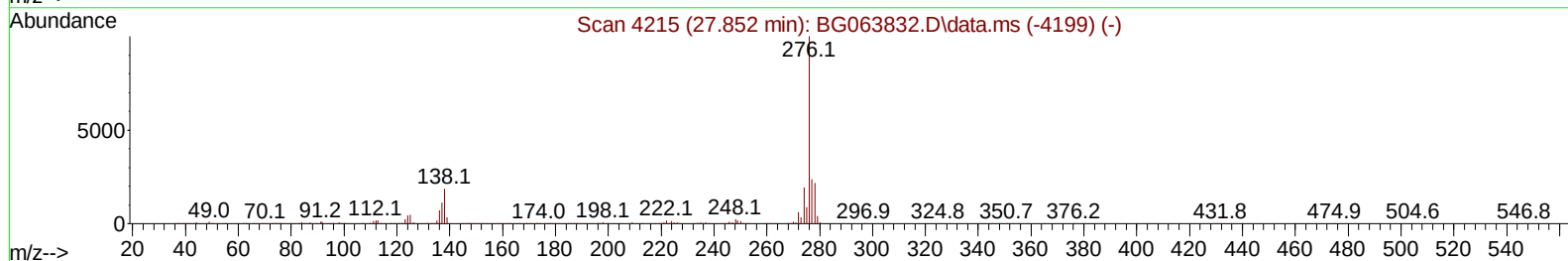
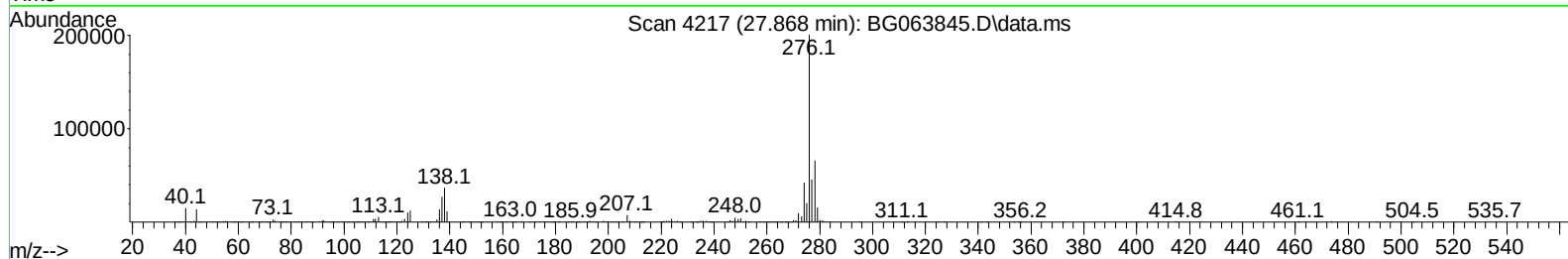
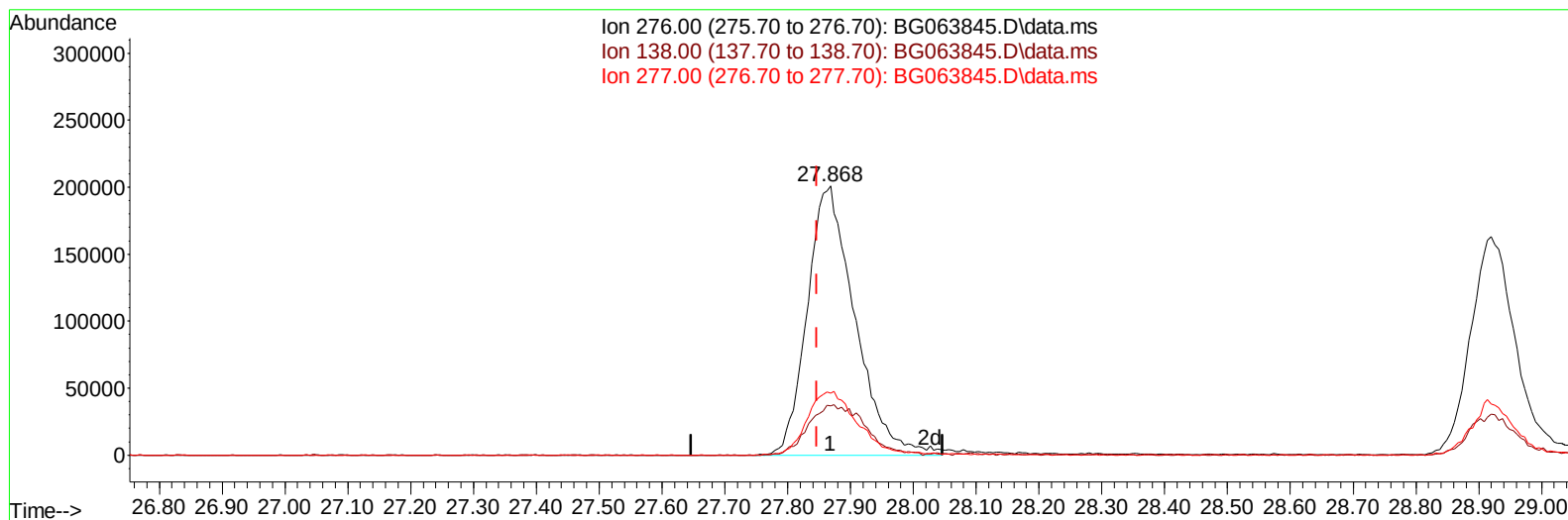
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TIC: BG063845.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

27.868min (+ 0.021) 21.72 ng/ul m

response 1058198

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	17.10	18.29
277.00	23.00	22.91
0.00	0.00	0.00

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Instrument :
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 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 27 02:45:20 2024
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 QLast Update : Tue Dec 24 00:44:39 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	88227	20.000	ng/ul	0.00
20) Naphthalene-d8	10.588	136	358736	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.437	164	277241	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	683480	20.000	ng/ul	0.00
79) Chrysene-d12	21.440	240	711419	20.000	ng/ul	0.00
88) Perylene-d12	24.460	264	747668	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	20385	9.054	ng/uL	0.00
4) Pyridine-d5	3.661	84	146719	20.658	ng/ul	0.00
7) Phenol-d5	6.981	99	159560	19.637	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	99837	17.790	ng/ul	0.00
11) 2-Chlorophenol-d4	7.333	132	102639	19.511	ng/ul	0.00
15) 4-Methylphenol-d8	8.514	113	119832	18.995	ng/ul	0.00
21) Nitrobenzene-d5	8.949	128	57056	22.275	ng/ul	0.00
24) 2-Nitrophenol-d4	9.678	143	64886	22.599	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.218	165	120554	21.687	ng/ul	0.00
31) 4-Chloroaniline-d4	10.724	131	167589	23.744	ng/ul	0.00
46) Dimethylphthalate-d6	13.843	166	395631	20.972	ng/ul	0.00
49) Acenaphthylene-d8	14.131	160	461387	21.455	ng/ul	0.00
54) 4-Nitrophenol-d4	14.678	143	55382	20.720	ng/ul	0.00
60) Fluorene-d10	15.436	176	355459	21.311	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.571	200	69981	19.720	ng/ul	0.00
73) Anthracene-d10	17.292	188	650174	22.126	ng/ul	0.00
81) Pyrene-d10	19.590	212	766435	20.703	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.255	264	745271	21.226	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.291	88	23540	8.835	ng/uL	87
5) Pyridine	3.685	79	164696	21.650	ng/ul	92
6) Benzaldehyde	6.940	77	97162	19.494	ng/ul	96
8) Phenol	7.010	94	162909	18.812	ng/ul	96
10) Bis(2-Chloroethyl)ether	7.222	93	125051	18.873	ng/ul	97
12) 2-Chlorophenol	7.369	128	106108	19.927	ng/ul	98
13) 2-Methylphenol	8.250	108	120627	19.272	ng/ul	98
14) 2,2'-oxybis(1-Chloropr...	8.315	45	183742	18.680	ng/ul	95
16) Acetophenone	8.614	105	206280	19.278	ng/ul	94
17) N-Nitroso-di-n-propyla...	8.597	70	116866	18.304	ng/ul	88
18) 4-Methylphenol	8.579	108	137813	20.118	ng/ul	97
19) Hexachloroethane	8.867	117	47140	18.017	ng/ul	91
22) Nitrobenzene	8.996	77	176988	20.692	ng/ul	98
23) Isophorone	9.513	82	324343	20.069	ng/ul	96
25) 2-Nitrophenol	9.707	139	66134	22.412	ng/ul	90
26) 2,4-Dimethylphenol	9.772	107	140504	20.984	ng/ul	91
27) Bis(2-Chloroethoxy)met...	9.995	93	181298	20.741	ng/ul	91
29) 2,4-Dichlorophenol	10.248	162	117419	21.258	ng/ul	94
30) Naphthalene	10.635	128	380028	21.188	ng/ul	99
32) 4-Chloroaniline	10.747	127	160087	23.838	ng/ul	99
33) Hexachlorobutadiene	10.917	225	84087	18.738	ng/ul	99
34) Caprolactam	11.505	113	44472m	23.292	ng/ul	
35) 4-Chloro-3-methylphenol	11.893	107	142521	21.978	ng/ul	95

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SSTDCCC020

Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 12/30/2024

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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.251	142	270481	20.983	ng/ul	99
37) 1-Methylnaphthalene	12.469	142	280268	21.345	ng/ul	100
39) 1,2,4,5-Tetrachloroben...	12.627	216	168986	19.040	ng/ul	98
40) Hexachlorocyclopentadiene	12.598	237	54890	15.645	ng/ul	94
41) 2,4,6-Trichlorophenol	12.874	196	101825	20.299	ng/ul	96
42) 2,4,5-Trichlorophenol	12.950	196	107969	20.624	ng/ul	99
43) 1,1'-Biphenyl	13.268	154	380498	20.811	ng/ul	97
44) 2-Chloronaphthalene	13.309	162	307356	20.934	ng/ul	96
45) 2-Nitroaniline	13.520	65	108386	22.039	ng/ul	94
47) Dimethylphthalate	13.890	163	392592	20.765	ng/ul	97
48) 2,6-Dinitrotoluene	14.020	165	79683	22.693	ng/ul	92
50) Acenaphthylene	14.161	152	491809	21.329	ng/ul	100
51) 3-Nitroaniline	14.349	138	56955	17.544	ng/ul	99
52) Acenaphthene	14.502	153	349869	21.407	ng/ul	98
53) 2,4-Dinitrophenol	14.572	184	29822	16.619	ng/ul	95
55) 4-Nitrophenol	14.684	109	57582	20.747	ng/ul	95
56) Dibenzofuran	14.842	168	500255	21.711	ng/ul	93
57) 2,4-Dinitrotoluene	14.813	165	120359	23.068	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.077	232	90951	20.155	ng/ul	97
59) Diethylphthalate	15.259	149	394218	21.568	ng/ul	97
61) Fluorene	15.489	166	408963	22.228	ng/ul	97
62) 4-Chlorophenyl-phenyle...	15.483	204	212511	20.732	ng/ul	94
63) 4-Nitroaniline	15.518	138	48340	14.873	ng/ul	94
66) 4,6-Dinitro-2-methylph...	15.588	198	72323	19.783	ng/ul#	95
67) N-Nitrosodiphenylamine	15.700	169	353939	21.385	ng/ul	98
68) 4-Bromophenyl-phenylether	16.382	248	140836	20.219	ng/ul	98
69) Hexachlorobenzene	16.505	284	160012	20.503	ng/ul	97
70) Atrazine	16.652	200	148849	21.275	ng/ul	94
71) Pentachlorophenol	16.858	266	46463	14.121	ng/ul	94
72) Phenanthrene	17.234	178	737155	22.381	ng/ul	97
74) Anthracene	17.328	178	748804	22.462	ng/ul	98
75) 1,2,3,4-Tetrachloroben...	13.232	216	169363	18.572	ng/ul	98
76) Pentachlorobenzene	14.766	250	180034	19.928	ng/ul	95
77) Carbazole	17.598	167	657317	24.359	ng/ul	98
78) Di-n-butylphthalate	18.156	149	736827	22.610	ng/ul	98
80) Fluoranthene	19.255	202	895994	21.294	ng/ul	98
82) Pyrene	19.619	202	955229	21.338	ng/ul#	94
83) Butylbenzylphthalate	20.512	149	316978	23.066	ng/ul	97
84) 3,3'-Dichlorobenzidine	21.346	252	235318	18.323	ng/ul	95
85) Benzo(a)anthracene	21.423	228	915682	20.538	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.335	149	459160	22.969	ng/ul	99
87) Chrysene	21.487	228	867009	20.403	ng/ul	97
89) Di-n-octyl phthalate	22.469	149	793831	24.599	ng/ul	100
90) Benzo(b)fluoranthene	23.509	252	914177	21.513	ng/ul	99
91) Benzo(k)fluoranthene	23.567	252	894133	21.094	ng/ul	98
93) Benzo(a)pyrene	24.319	252	830270	20.816	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	27.868	276	1058198m	21.724	ng/ul	
95) Dibenzo(a,h)anthracene	27.909	278	851475	21.846	ng/ul	96
96) Benzo(g,h,i)perylene	28.920	276	812619	21.023	ng/ul	97

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

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