

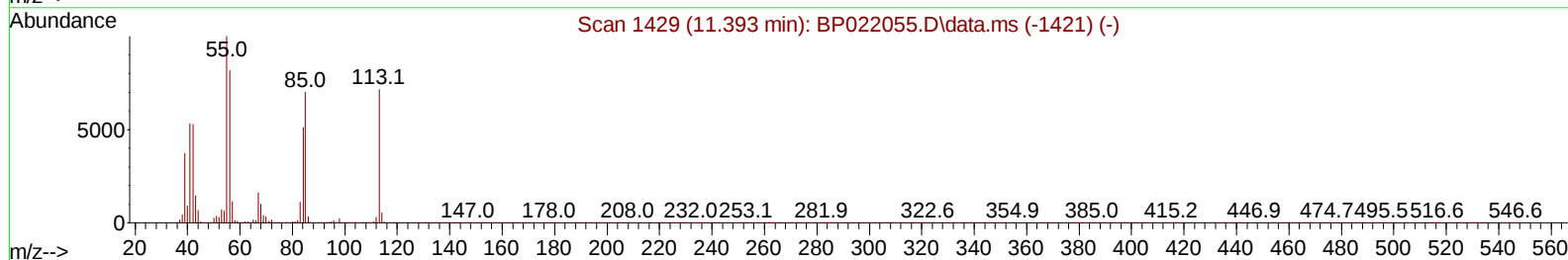
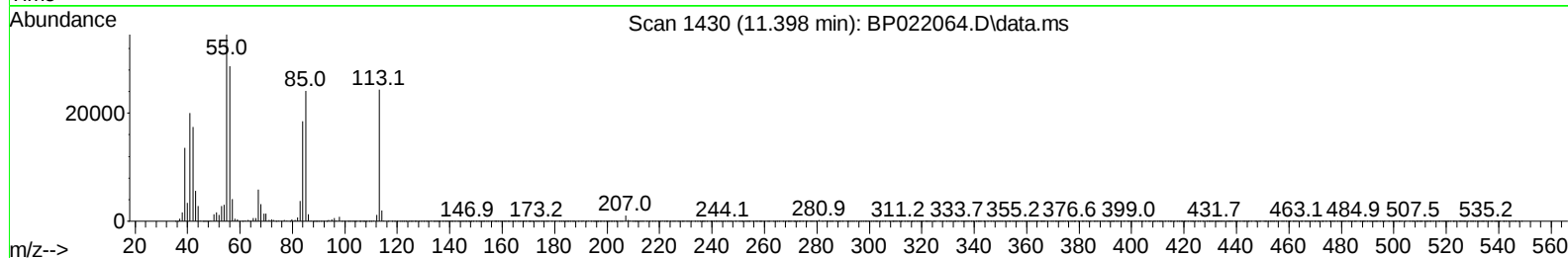
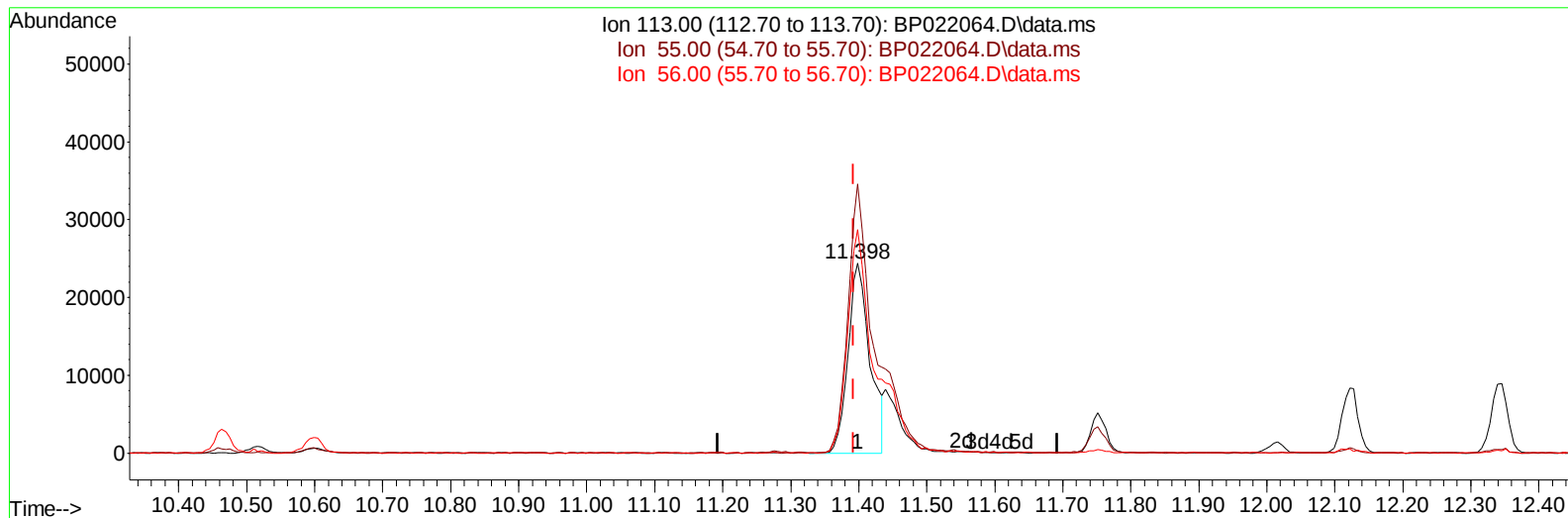
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022064.D
Acq On : 26 Sep 2024 15:51
Operator : RC/JU
Sample : PB163588BS
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SLCS588

Manual IntegrationsAPPROVED

Quant Time: Sep 27 01:17:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2024
Supervised By :mohammad ahmed 09/28/2024



TIC: BP022064.D\data.ms

(34) Caprolactam

11.398min (+ 0.006) 24.95 ng/ul

response 54543

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	136.10	141.83
56.00	127.30	117.86
0.00	0.00	0.00

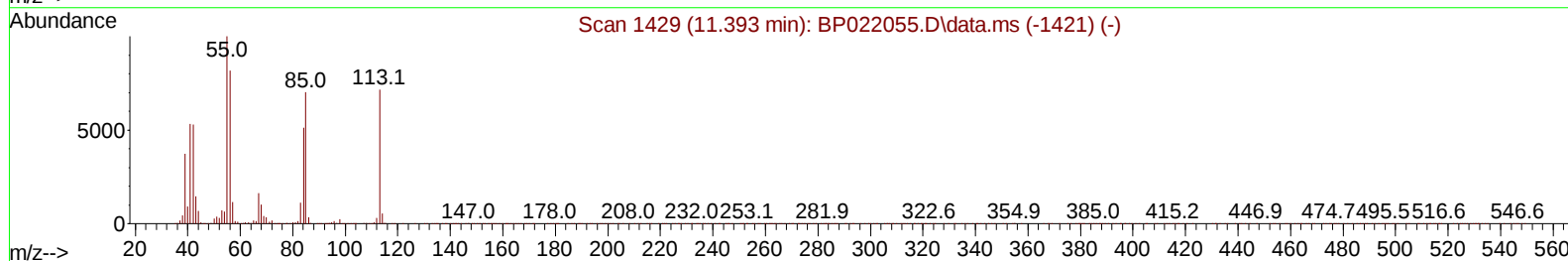
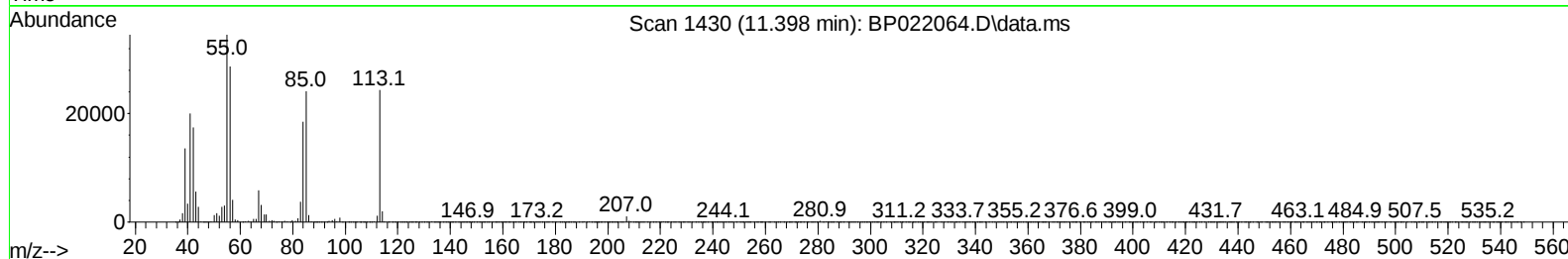
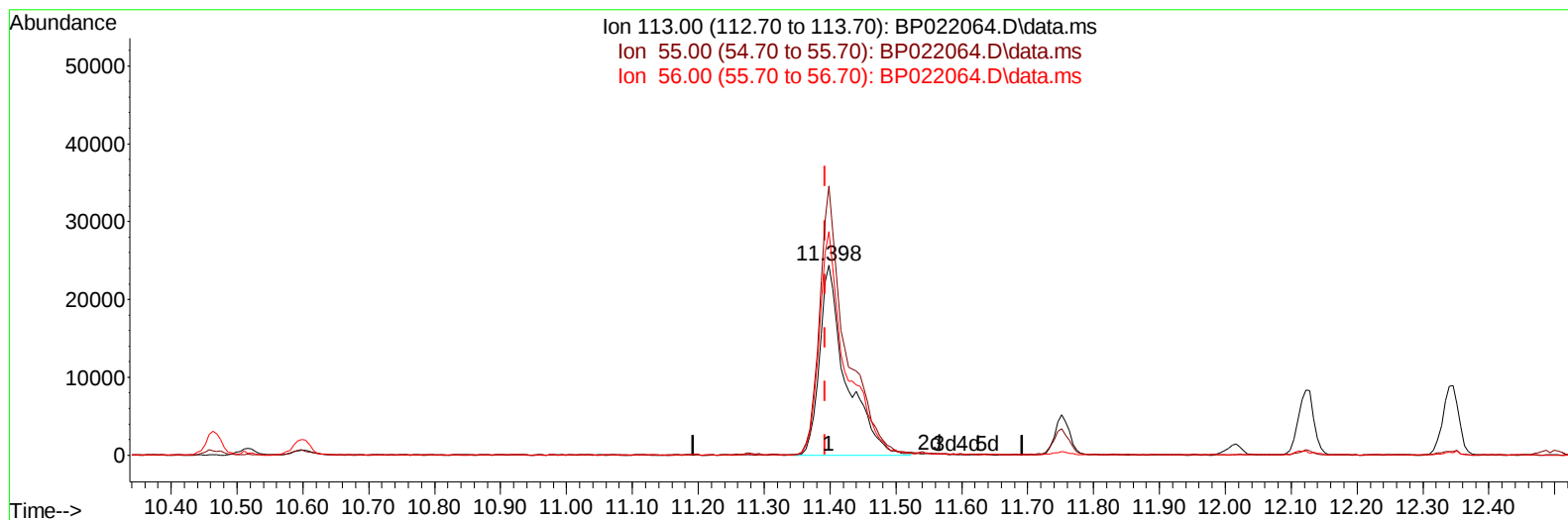
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TIC: BP022064.D\data.ms

(34) Caprolactam

11.398min (+ 0.006) 31.24 ng/ul m

response 68300

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	136.10	141.83
56.00	127.30	117.86
0.00	0.00	0.00

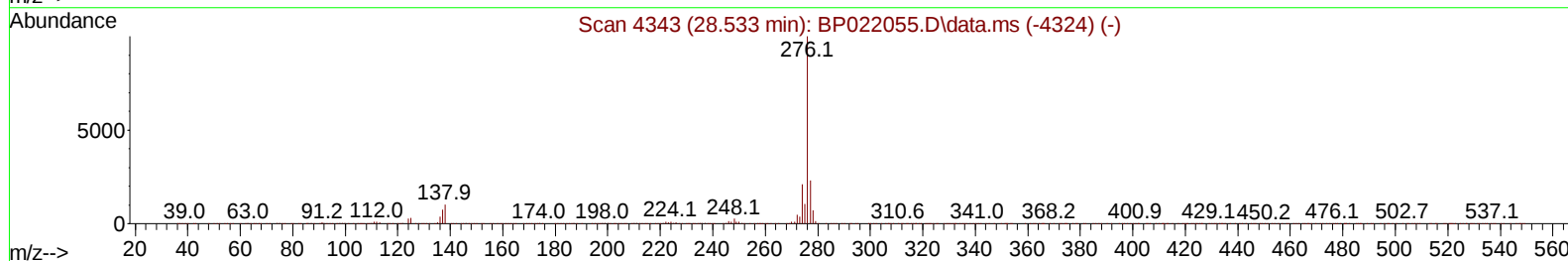
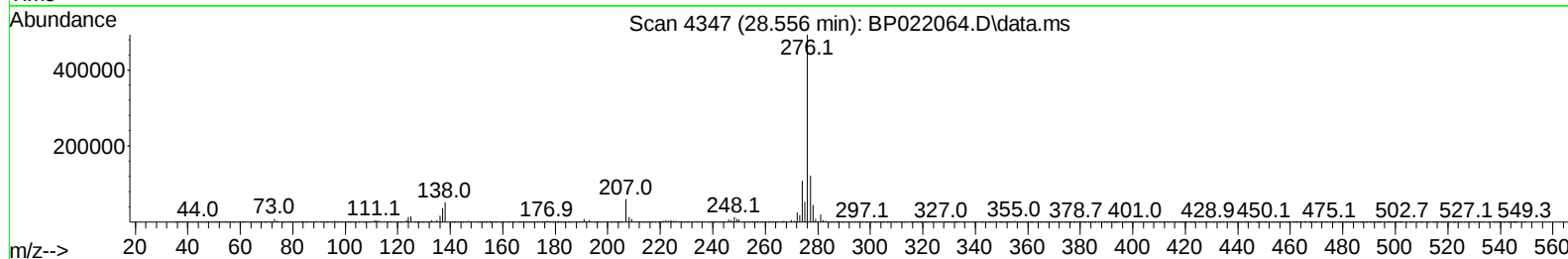
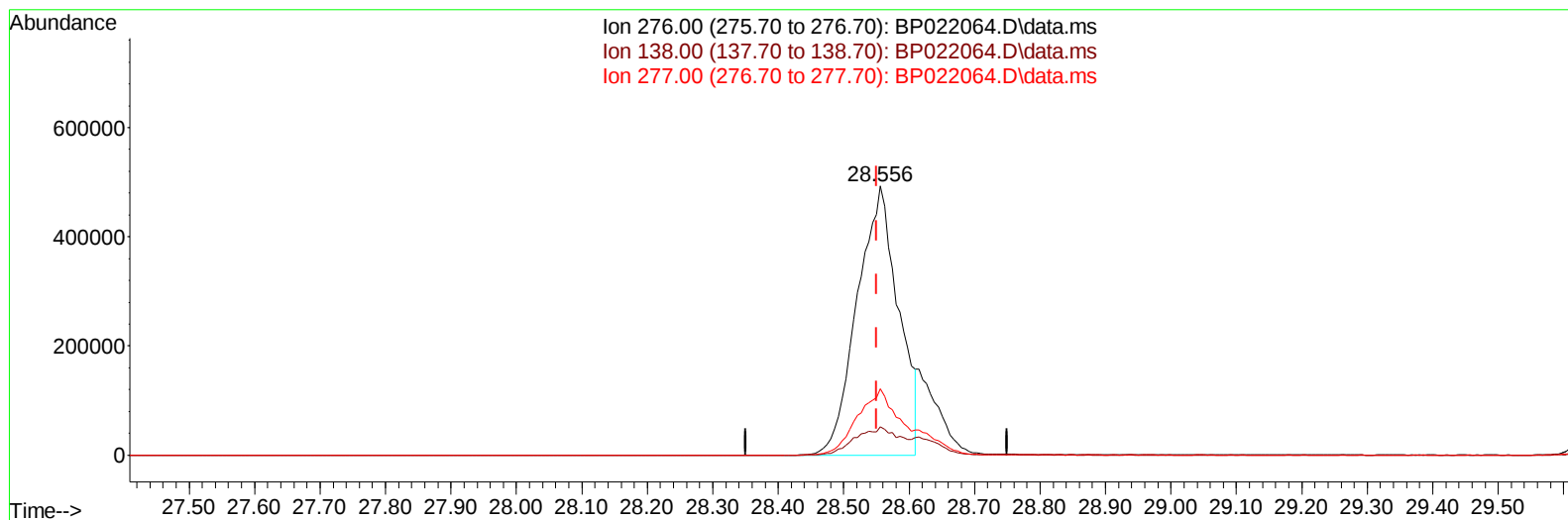
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Operator : RC/JU
Sample : PB163588BS
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP022064.D\data.ms

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

28.556min (+ 0.006) 24.69 ng/u1

response 2148803

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	11.50	10.41
277.00	24.20	24.62
0.00	0.00	0.00

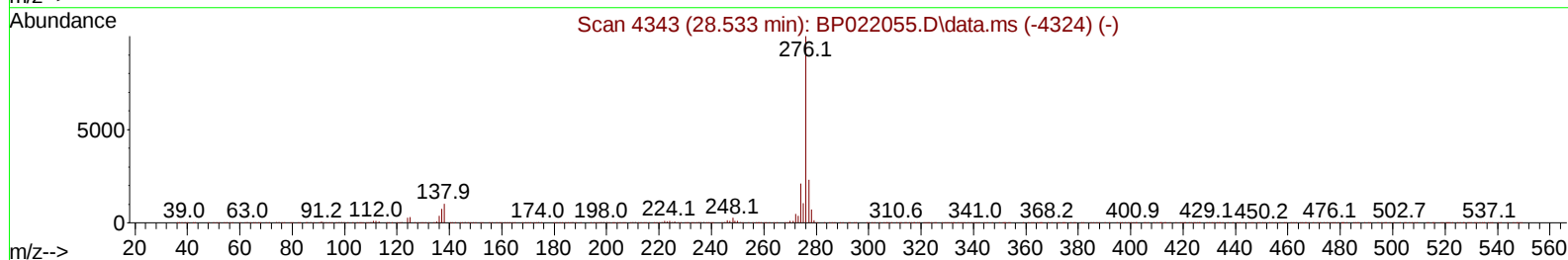
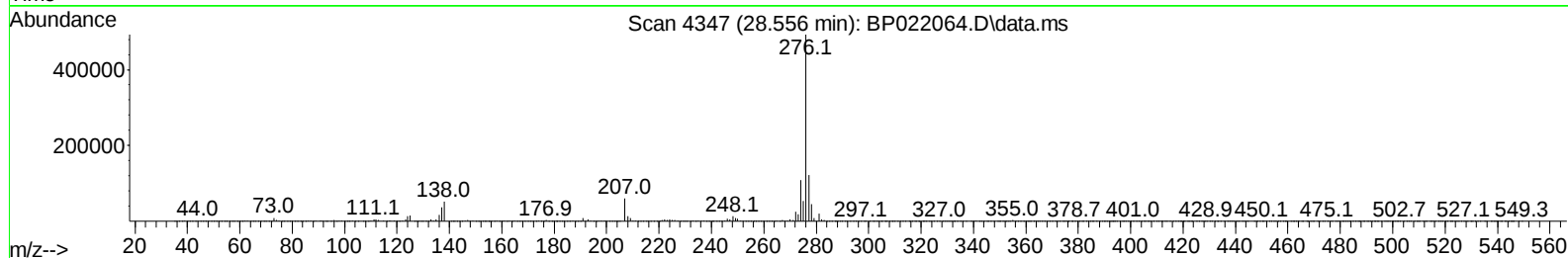
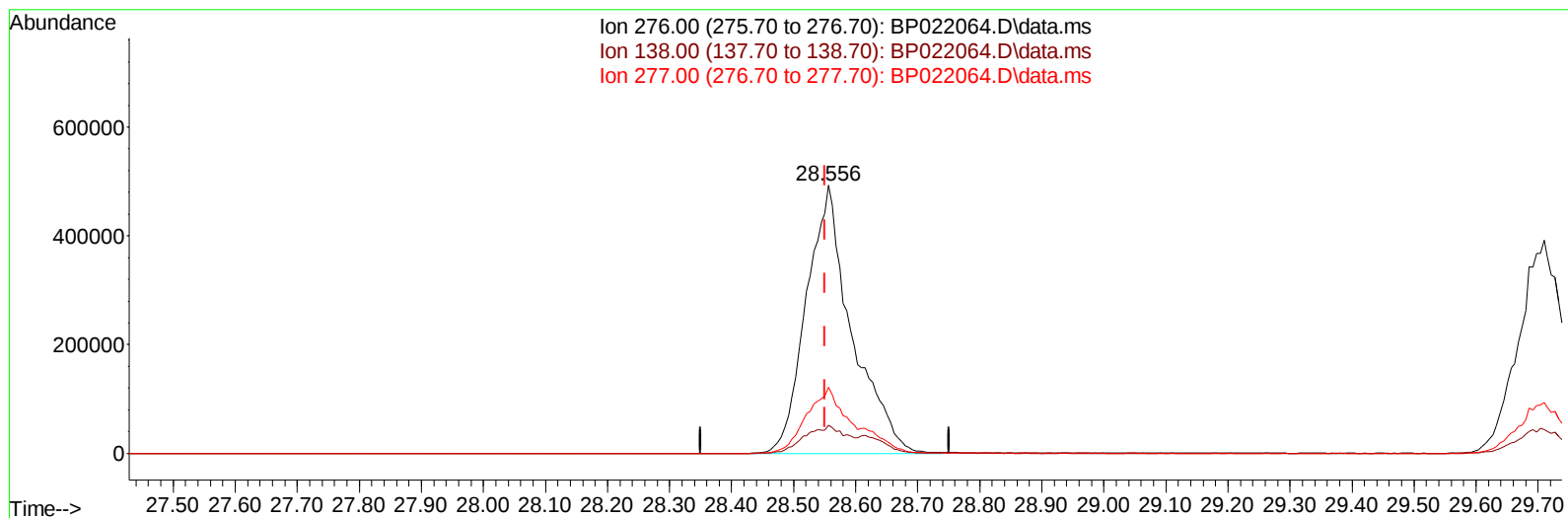
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
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Operator : RC/JU
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Misc :
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Instrument :
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Supervised By :mohammad ahmed 09/28/2024



TIC: BP022064.D\data.ms

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

28.556min (+ 0.006) 28.71 ng/ul m

response 2498272

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	11.50	10.41
277.00	24.20	24.62
0.00	0.00	0.00

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 Operator : RC/JU
 Sample : PB163588BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLCS588

Manual IntegrationsAPPROVED

Quant Time: Sep 27 01:34:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2024
 Supervised By :mohammad ahmed 09/28/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	99818	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	393605	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	328000	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	851903	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1030931	20.000	ng/ul	0.00
88) Perylene-d12	24.827	264	1166699	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	12998	5.101	ng/uL	0.00
4) Pyridine-d5	3.634	84	180228	24.742	ng/ul	0.00
7) Phenol-d5	6.893	99	230327	28.208	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	138983	27.643	ng/ul	0.00
11) 2-Chlorophenol-d4	7.252	132	177818	30.100	ng/ul	0.00
15) 4-Methylphenol-d8	8.410	113	185562	28.680	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	88161	29.926	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	103248	30.910	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	221894	30.589	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	242049	27.966	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	754737	29.767	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	799709	29.791	ng/ul	0.00
54) 4-Nitrophenol-d4	14.534	143	123910	28.571	ng/ul	0.00
60) Fluorene-d10	15.322	176	680753	30.353	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	138245	26.138	ng/ul	0.00
73) Anthracene-d10	17.222	188	1187435	29.857	ng/ul	0.00
81) Pyrene-d10	19.586	212	1610443	30.152	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	1895567	30.890	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	27182	9.847	ng/uL	99
5) Pyridine	3.652	79	175273	23.684	ng/ul	98
6) Benzaldehyde	6.858	77	150862	32.661	ng/ul	99
8) Phenol	6.916	94	236472	27.751	ng/ul	96
10) Bis(2-Chloroethyl)ether	7.134	93	177272	27.181	ng/ul	97
12) 2-Chlorophenol	7.281	128	177908	29.049	ng/ul	99
13) 2-Methylphenol	8.146	108	172441	27.903	ng/ul	95
14) 2,2'-oxybis(1-Chloropr...	8.216	45	156488	25.865	ng/ul	95
16) Acetophenone	8.516	105	306026	28.435	ng/ul	98
17) N-Nitroso-di-n-propyla...	8.505	70	160599	28.784	ng/ul	95
18) 4-Methylphenol	8.469	108	190246	28.229	ng/ul	99
19) Hexachloroethane	8.769	117	81901	28.329	ng/ul	92
22) Nitrobenzene	8.887	77	253442	28.138	ng/ul	100
23) Isophorone	9.410	82	440426	28.230	ng/ul	98
25) 2-Nitrophenol	9.587	139	107359	30.039	ng/ul	94
26) 2,4-Dimethylphenol	9.652	107	197379	24.379	ng/ul	98
27) Bis(2-Chloroethoxy)met...	9.881	93	248262	27.606	ng/ul	98
29) 2,4-Dichlorophenol	10.122	162	215975	29.968	ng/ul	99
30) Naphthalene	10.516	128	576926	27.857	ng/ul	99
32) 4-Chloroaniline	10.622	127	218214	25.456	ng/ul	99
33) Hexachlorobutadiene	10.799	225	182964	28.594	ng/ul	96
34) Caprolactam	11.398	113	68300m	31.241	ng/ul	
35) 4-Chloro-3-methylphenol	11.751	107	233745	29.979	ng/ul	99

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Manual IntegrationsAPPROVED

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Quant Time: Sep 27 01:34:00 2024
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.122	142	445664	28.956	ng/ul	98
37) 1-Methylnaphthalene	12.345	142	447977	28.860	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.493	216	337050	27.716	ng/ul	98
40) Hexachlorocyclopentadiene	12.475	237	185907	24.304	ng/ul	97
41) 2,4,6-Trichlorophenol	12.740	196	211270	29.061	ng/ul	98
42) 2,4,5-Trichlorophenol	12.810	196	234040	29.729	ng/ul	98
43) 1,1'-Biphenyl	13.140	154	647316	27.598	ng/ul	98
44) 2-Chloronaphthalene	13.181	162	528218	27.970	ng/ul	99
45) 2-Nitroaniline	13.387	65	168631	30.881	ng/ul	98
47) Dimethylphthalate	13.769	163	723782	28.421	ng/ul	99
48) 2,6-Dinitrotoluene	13.887	165	149083	30.537	ng/ul	96
50) Acenaphthylene	14.034	152	791322	28.267	ng/ul	99
51) 3-Nitroaniline	14.222	138	131960	29.689	ng/ul	96
52) Acenaphthene	14.381	153	573798	28.210	ng/ul	97
53) 2,4-Dinitrophenol	14.434	184	59721	17.978	ng/ul	94
55) 4-Nitrophenol	14.545	109	151217	30.819	ng/ul	89
56) Dibenzofuran	14.716	168	860115	28.442	ng/ul	100
57) 2,4-Dinitrotoluene	14.681	165	234915	31.224	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	14.951	232	231311	30.525	ng/ul	99
59) Diethylphthalate	15.163	149	746758	29.629	ng/ul	98
61) Fluorene	15.381	166	734470	29.205	ng/ul	98
62) 4-Chlorophenyl-phenyle...	15.375	204	432226	29.852	ng/ul	98
63) 4-Nitroaniline	15.392	138	150115	33.183	ng/ul	94
66) 4,6-Dinitro-2-methylph...	15.469	198	143733	24.962	ng/ul#	99
67) N-Nitrosodiphenylamine	15.592	169	640954	28.566	ng/ul	99
68) 4-Bromophenyl-phenylether	16.292	248	304786	29.792	ng/ul	99
69) Hexachlorobenzene	16.404	284	375534	30.231	ng/ul	98
70) Atrazine	16.575	200	260145	25.990	ng/ul	99
71) Pentachlorophenol	16.757	266	209202	25.242	ng/ul	99
72) Phenanthrene	17.163	178	1299998	29.104	ng/ul	99
74) Anthracene	17.257	178	1325034	29.151	ng/ul	100
75) 1,2,3,4-Tetrachloroben...	13.104	216	338011	26.441	ng/uL	98
76) Pentachlorobenzene	14.639	250	381677	27.150	ng/uL	99
77) Carbazole	17.533	167	1178822	29.425	ng/ul	100
78) Di-n-butylphthalate	18.122	149	1420953	30.765	ng/ul	99
80) Fluoranthene	19.239	202	1788887	29.304	ng/ul	100
82) Pyrene	19.622	202	1926695	29.321	ng/ul	96
83) Butylbenzylphthalate	20.569	149	710884	31.042	ng/ul	99
84) 3,3'-Dichlorobenzidine	21.451	252	628537	26.628	ng/ul	98
85) Benzo(a)anthracene	21.527	228	2069839	29.133	ng/ul	100
86) Bis(2-ethylhexyl)phtha...	21.463	149	1039206	30.929	ng/ul	100
87) Chrysene	21.592	228	1969232	29.533	ng/ul	100
89) Di-n-octyl phthalate	22.710	149	1822902	31.628	ng/ul	100
90) Benzo(b)fluoranthene	23.780	252	2139160	29.536	ng/ul	99
91) Benzo(k)fluoranthene	23.857	252	2156579	29.876	ng/ul	99
93) Benzo(a)pyrene	24.674	252	1983877	29.727	ng/ul	100
94) Indeno(1,2,3-cd)pyrene	28.556	276	2498272m	28.711	ng/ul	
95) Dibenzo(a,h)anthracene	28.615	278	2004265	28.593	ng/ul	98
96) Benzo(g,h,i)perylene	29.709	276	1928195	28.558	ng/ul	100

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

Instrument :

BNA_P

ClientSampleId :

SLCS588

Manual IntegrationsAPPROVED

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