

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018538.D
 Acq On : 03 Dec 2023 04:30
 Operator : MA/JU
 Sample : 05411-16
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QK1

Quant Time: Dec 03 23:49:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	404566	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.657	136	1641500	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	968155	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2020619	20.000	ng/u1	0.00
79) Chrysene-d12	21.315	240	1953665	20.000	ng/u1	-0.01
88) Perylene-d12	23.692	264	2319002	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	45642	3.844	ng/uL	0.00
4) Pyridine-d5	3.693	84	295193	9.283	ng/u1	0.00
7) Phenol-d5	7.022	99	197055	4.854	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.193	67	627622	26.082	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.387	132	574528	18.135	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	368863	11.226	ng/u1	-0.01
21) Nitrobenzene-d5	9.022	128	367624	25.163	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	338520	22.349	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	645693	20.847	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.792	131	912810	22.551	ng/u1	-0.01
46) Dimethylphthalate-d6	13.904	166	2208536	25.559	ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	2499446	26.269	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	50276	3.545	ng/u1	-0.01
60) Fluorene-d10	15.480	176	1924798	26.537	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	219861	18.147	ng/u1	-0.01
73) Anthracene-d10	17.333	188	2970778	27.966	ng/u1	0.00
81) Pyrene-d10	19.568	212	3491107	22.956	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	3757242	27.787	ng/u1	-0.01

Target Compounds	Qvalue

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

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