

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018529.D
 Acq On : 02 Dec 2023 22:50
 Operator : MA/JU
 Sample : 05411-10
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QH5

Quant Time: Dec 02 23:25:46 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	7.858	152	413849	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.658	136	1830885	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.487	164	1194548	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2364014	20.000	ng/u1	0.00
79) Chrysene-d12	21.316	240	1687444	20.000	ng/u1	-0.01
88) Perylene-d12	23.692	264	2017466	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	46945	3.866	ng/uL	0.00
4) Pyridine-d5	3.693	84	277780	8.539	ng/u1	0.00
7) Phenol-d5	7.022	99	237191	5.711	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.193	67	728394	29.591	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.387	132	666814	20.576	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	457814	13.621	ng/u1	-0.01
21) Nitrobenzene-d5	9.022	128	471563	28.938	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	470746	27.864	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	835592	24.188	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.799	131	1181463	26.169	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	3200038	30.015	ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	3531734	30.084	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	66499	3.800	ng/u1	-0.01
60) Fluorene-d10	15.481	176	2667159	29.803	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	331373	23.378	ng/u1	-0.01
73) Anthracene-d10	17.334	188	3875887	31.187	ng/u1	0.00
81) Pyrene-d10	19.569	212	3915555	29.809	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	3726732	31.681	ng/u1	-0.01

Target Compounds Qvalue

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

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