

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120123\
 Data File : BP018507.D
 Acq On : 02 Dec 2023 09:49
 Operator : MA/JU
 Sample : 05328-02
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49P9

Quant Time: Dec 02 21:29:29 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	7.863	152	515383	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	2113116	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	1277160	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	2689156	20.000	ng/ul	0.00
79) Chrysene-d12	21.321	240	2690155	20.000	ng/ul	0.00
88) Perylene-d12	23.698	264	3217074	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	52689	3.484	ng/uL	0.00
4) Pyridine-d5	3.693	84	177673	4.386	ng/ul	0.00
7) Phenol-d5	7.028	99	197997	3.828	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	656908	21.430	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	602543	14.930	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	370648	8.855	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	396270	21.070	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	385159	19.753	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	732947	18.383	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	760831	14.601	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	2587103	22.696	ng/ul	0.00
49) Acenaphthylene-d8	14.181	160	2818965	22.459	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	65625	3.507	ng/ul	-0.01
60) Fluorene-d10	15.481	176	2226999	23.275	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	310711	19.270	ng/ul	0.00
73) Anthracene-d10	17.333	188	3418383	24.180	ng/ul	0.00
81) Pyrene-d10	19.569	212	4403785	21.030	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.545	264	4775717	25.460	ng/ul	0.00

Target Compounds Qvalue

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

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