

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018287.D
 Acq On : 22 Nov 2023 22:02
 Operator : MA/JU
 Sample : 05226-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG47

Quant Time: Nov 22 23:31:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	368003	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1517702	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	916537	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1665533	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1130040	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1353796	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	39051	3.975	ng/uL	0.00
4) Pyridine-d5	3.687	84	278218	10.304	ng/u1	0.00
7) Phenol-d5	7.034	99	256368	7.623	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	707213	35.165	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	681701	24.142	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	450589	16.602	ng/u1	0.00
21) Nitrobenzene-d5	9.034	128	447003	36.261	ng/u1	0.00
24) 2-Nitrophenol-d4	9.758	143	398843	35.033	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	775549	28.236	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	1162269	31.375	ng/u1	0.00
46) Dimethylphthalate-d6	13.928	166	2714066	33.279	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	3159440	34.000	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	55320	4.914	ng/u1	-0.01
60) Fluorene-d10	15.498	176	2285596	33.507	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	188790	26.627	ng/u1	0.00
73) Anthracene-d10	17.357	188	3287647	36.931	ng/u1	0.00
81) Pyrene-d10	19.592	212	3186006	35.425	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	3034576	38.663	ng/u1	0.00

Target Compounds	Qvalue

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

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