

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018276.D
 Acq On : 22 Nov 2023 14:55
 Operator : MA/JU
 Sample : PB157005BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK005

Quant Time: Nov 22 23:34:04 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	7.869	152	322164	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.675	136	1349707	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	886626	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.263	188	1754871	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	1029257	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	1144057	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	51871	6.030	ng/uL	0.00
4) Pyridine-d5	3.681	84	703092	29.744	ng/u1	0.00
7) Phenol-d5	7.028	99	853114	28.975	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	528671	30.028	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	695939	28.152	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	694143	29.214	ng/u1	-0.01
21) Nitrobenzene-d5	9.028	128	336460	30.691	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.752	143	322432	31.847	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	675587	27.658	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.810	131	1056465	32.068	ng/u1	0.00
46) Dimethylphthalate-d6	13.928	166	2344167	29.713	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	2638043	29.346	ng/u1	0.00
54) 4-Nitrophenol-d4	14.693	143	269332	24.729	ng/u1	0.00
60) Fluorene-d10	15.504	176	1920927	29.111	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	153004	20.481	ng/u1	0.00
73) Anthracene-d10	17.363	188	2780580	29.645	ng/u1	0.00
81) Pyrene-d10	19.604	212	2590293	31.621	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	1987428	29.964	ng/u1	0.00

Target Compounds Qvalue

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

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