

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018321.D
 Acq On : 24 Nov 2023 09:42
 Operator : MA/JU
 Sample : PB157173BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK173

Quant Time: Nov 24 21:54:11 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	384859	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.669	136	1589770	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.504	164	1022862	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	2028616	20.000	ng/ul	-0.01
79) Chrysene-d12	21.357	240	1217001	20.000	ng/ul	0.00
88) Perylene-d12	23.786	264	1273969	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	59276	5.769	ng/uL	0.00
4) Pyridine-d5	3.687	84	721957	25.567	ng/ul	0.00
7) Phenol-d5	7.028	99	1001130	28.464	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	612460	29.120	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	823796	27.896	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	811419	28.587	ng/ul	-0.01
21) Nitrobenzene-d5	9.028	128	392604	30.405	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.752	143	371937	31.189	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	793605	27.584	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.805	131	950073	24.484	ng/ul	-0.01
46) Dimethylphthalate-d6	13.922	166	2754636	30.266	ng/ul	-0.01
49) Acenaphthylene-d8	14.199	160	3018748	29.109	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	304608	24.243	ng/ul	-0.01
60) Fluorene-d10	15.498	176	2262173	29.716	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	205760	23.827	ng/ul	-0.01
73) Anthracene-d10	17.351	188	3302423	30.457	ng/ul	-0.01
81) Pyrene-d10	19.592	212	3190563	32.940	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	2302948	31.180	ng/ul	0.00

Target Compounds Qvalue

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

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