

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121323\
 Data File : BP018846.D
 Acq On : 14 Dec 2023 17:01
 Operator : MA/JU
 Sample : PB157722BL
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK722

Quant Time: Dec 14 21:55:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 13 23:34:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	237800	20.000	ng/ul	0.00
20) Naphthalene-d8	10.628	136	827399	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	464049	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1042165	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1186966	20.000	ng/ul	0.00
88) Perylene-d12	23.662	264	1425861	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	34815	4.989	ng/uL	0.00
4) Pyridine-d5	3.675	84	433034	23.167	ng/ul	0.00
7) Phenol-d5	7.005	99	513007	21.497	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.169	67	300153	21.221	ng/ul	0.00
11) 2-Chlorophenol-d4	7.363	132	410969	22.070	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	317406	16.435	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	128572	17.459	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	146192	19.148	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	354669	22.718	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	474644	23.264	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1000561	24.158	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1100884	24.139	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	159359	23.441	ng/ul	0.00
60) Fluorene-d10	15.451	176	867092	24.941	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	116571	18.655	ng/ul	0.00
73) Anthracene-d10	17.310	188	1501842	27.412	ng/ul	0.00
81) Pyrene-d10	19.545	212	1993639	21.577	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	2296295	27.620	ng/ul	-0.01

Target Compounds Qvalue

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

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