

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018078.D
 Acq On : 11 Nov 2023 22:33
 Operator : MA/JU
 Sample : PB156719BL
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK719

Quant Time: Nov 12 08:47:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	70406	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	282933	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	178605	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.304	188	403980	20.000	ng/u1	0.00
79) Chrysene-d12	21.427	240	399563	20.000	ng/u1	0.00
88) Perylene-d12	23.933	264	463772	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	10233	5.127	ng/uL	0.00
4) Pyridine-d5	3.664	84	129448	24.875	ng/u1	0.00
7) Phenol-d5	7.016	99	152458	22.430	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	99108	24.811	ng/u1	0.00
11) 2-Chlorophenol-d4	7.369	132	127467	23.681	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	127307	22.856	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	61517	24.133	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	71035	24.598	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	115566	22.453	ng/u1	0.00
31) 4-Chloroaniline-d4	10.851	131	160664	22.762	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	402917	24.450	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	427389	23.811	ng/u1	0.00
54) 4-Nitrophenol-d4	14.798	143	43517	17.672	ng/u1	0.02
60) Fluorene-d10	15.522	176	329093	24.188	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	58363	19.984	ng/u1	0.00
73) Anthracene-d10	17.404	188	544989	24.944	ng/u1	0.00
81) Pyrene-d10	19.657	212	668496	25.047	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.762	264	690140	25.592	ng/u1	-0.01

Target Compounds Qvalue

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

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