

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017175.D
 Acq On : 14 Sep 2023 13:59
 Operator : MA/JU
 Sample : PB155433BL
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK433

Quant Time: Sep 14 22:47:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	199556	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	826139	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.610	164	494082	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.381	188	1050018	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	833630	20.000	ng/ul	0.00
88) Perylene-d12	24.010	264	917719	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	30776	6.481	ng/uL	0.00
4) Pyridine-d5	3.746	84	296675	21.839	ng/ul	0.00
7) Phenol-d5	7.105	99	529340	32.622	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	334605	32.992	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	419097	32.376	ng/ul	0.00
15) 4-Methylphenol-d8	8.664	113	410823	30.549	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	196176	33.912	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	216453	35.049	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	390599	32.704	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	377689	20.980	ng/ul	0.00
46) Dimethylphthalate-d6	14.028	166	1241964	35.323	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	1352014	33.889	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	163496	24.060	ng/ul	0.00
60) Fluorene-d10	15.604	176	1054248	34.678	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.734	200	135435	23.086	ng/ul	-0.01
73) Anthracene-d10	17.481	188	1653730	37.068	ng/ul	0.00
81) Pyrene-d10	19.722	212	1810156	41.404	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	1650640	38.108	ng/ul	0.00

Target Compounds Qvalue

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

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