

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017151.D
 Acq On : 13 Sep 2023 16:16
 Operator : MA/JU
 Sample : PB155421BL
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK421

Quant Time: Sep 13 23:57:49 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	168901	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	716746	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	455926	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	1072489	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1064324	20.000	ng/u1	0.00
88) Perylene-d12	24.010	264	1174859	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	25146	6.256	ng/uL	0.00
4) Pyridine-d5	3.746	84	248119	21.580	ng/u1	0.00
7) Phenol-d5	7.105	99	415956	30.287	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	262031	30.525	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	325046	29.668	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	325785	28.622	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	154504	30.785	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	163611	30.536	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	305458	29.478	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	313139	20.049	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1051455	32.408	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1125628	30.576	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	160656	25.620	ng/u1	0.00
60) Fluorene-d10	15.610	176	918452	32.739	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	129326	21.583	ng/u1	-0.01
73) Anthracene-d10	17.481	188	1550565	34.028	ng/u1	0.00
81) Pyrene-d10	19.722	212	1919570	34.390	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	1896511	34.201	ng/u1	-0.01

Target Compounds Qvalue

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

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