

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017168.D
 Acq On : 14 Sep 2023 09:59
 Operator : MA/JU
 Sample : PB155343BL
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK343

Quant Time: Sep 14 22:54:18 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	7.940	152	197671	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.769	136	848783	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.610	164	562198	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	1326536	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	1229408	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	1261138	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	31511	6.699	ng/uL	-0.01
4) Pyridine-d5	3.740	84	304376	22.620	ng/u1	-0.01
7) Phenol-d5	7.087	99	529346	32.933	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.281	67	336731	33.518	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.457	132	421906	32.903	ng/u1	-0.02
15) 4-Methylphenol-d8	8.652	113	428593	32.174	ng/u1	-0.02
21) Nitrobenzene-d5	9.140	128	204787	34.456	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.857	143	224790	35.427	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	406738	33.146	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.934	131	407079	22.009	ng/u1	-0.01
46) Dimethylphthalate-d6	14.028	166	1456422	36.404	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1544566	34.025	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	221936	28.703	ng/u1	0.00
60) Fluorene-d10	15.610	176	1236941	35.758	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	173555	23.417	ng/u1	0.00
73) Anthracene-d10	17.486	188	2081122	36.925	ng/u1	0.00
81) Pyrene-d10	19.727	212	2493741	38.677	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	2226981	37.413	ng/u1	0.00

Target Compounds Qvalue

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

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