

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
 Data File : BP017239.D
 Acq On : 20 Sep 2023 15:09
 Operator : MA/JU
 Sample : PB155579BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK579

Quant Time: Sep 20 22:30:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 20 14:34:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	257055	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	1081165	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.604	164	688379	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	1577032	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1566832	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1798790	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	31867	5.092	ng/uL	0.00
4) Pyridine-d5	3.740	84	466045	26.638	ng/u1	0.00
7) Phenol-d5	7.093	99	583352	27.631	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	360438	28.290	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	471995	28.188	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	466653	28.431	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	228992	29.297	ng/u1	0.00
24) 2-Nitrophenol-d4	9.852	143	256049	30.170	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	444456	27.719	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	678016	29.191	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	1501282	30.444	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	1617160	28.513	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	250842	29.197	ng/u1	0.00
60) Fluorene-d10	15.598	176	1273234	29.991	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	227717	25.619	ng/u1	0.00
73) Anthracene-d10	17.475	188	2062148	29.483	ng/u1	0.00
81) Pyrene-d10	19.716	212	2594803	30.624	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2697025	31.141	ng/u1	0.00

Target Compounds	Qvalue

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

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