

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016869.D
 Acq On : 30 Aug 2023 13:13
 Operator : MA/JU
 Sample : PB155122BL
 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK122

Quant Time: Aug 31 01:20:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 30 02:25:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	349131	20.000	ng/u1	0.00
20) Naphthalene-d8	10.852	136	1532429	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.675	164	951037	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	2068779	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	1528265	20.000	ng/u1	0.00
88) Perylene-d12	24.110	264	1396685	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	61562	7.288	ng/uL	0.00
4) Pyridine-d5	3.811	84	846952	32.809	ng/u1	0.00
7) Phenol-d5	7.164	99	1097853	35.363	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	697298	35.241	ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	828076	35.476	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	867973	34.249	ng/u1	0.00
21) Nitrobenzene-d5	9.222	128	417068	35.996	ng/u1	0.00
24) 2-Nitrophenol-d4	9.934	143	449490	36.804	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.458	165	806712	35.430	ng/u1	0.00
31) 4-Chloroaniline-d4	11.010	131	1150100	32.938	ng/u1	0.00
46) Dimethylphthalate-d6	14.087	166	2576979	35.930	ng/u1	0.00
49) Acenaphthylene-d8	14.375	160	2935900	36.211	ng/u1	0.00
54) 4-Nitrophenol-d4	14.881	143	375440	25.326	ng/u1	0.00
60) Fluorene-d10	15.669	176	2251558	37.279	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	276329	22.429	ng/u1	0.00
73) Anthracene-d10	17.545	188	3536724	38.569	ng/u1	0.00
81) Pyrene-d10	19.781	212	3844057	45.798	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.939	264	2688592	38.430	ng/u1	0.00

Target Compounds Qvalue

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

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