

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
 Data File : BP017646.D
 Acq On : 11 Oct 2023 14:39
 Operator : MA/JU
 Sample : PB156277BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK277

Quant Time: Oct 12 02:34:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 04:45:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	105429	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	425306	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.634	164	276346	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.428	188	659604	20.000	ng/u1	0.01
79) Chrysene-d12	21.569	240	669150	20.000	ng/u1	0.02
88) Perylene-d12	24.157	264	729282	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	17954	6.909	ng/uL	0.00
4) Pyridine-d5	3.723	84	230397	31.495	ng/u1	0.00
7) Phenol-d5	7.093	99	273641	30.692	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	180220	32.593	ng/u1	0.01
11) 2-Chlorophenol-d4	7.458	132	228324	32.556	ng/u1	0.01
15) 4-Methylphenol-d8	8.658	113	221020	30.776	ng/u1	0.01
21) Nitrobenzene-d5	9.146	128	104856	31.996	ng/u1	0.01
24) 2-Nitrophenol-d4	9.863	143	115486	32.086	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.393	165	214932	30.746	ng/u1	0.01
31) 4-Chloroaniline-d4	10.946	131	299844	30.116	ng/u1	0.01
46) Dimethylphthalate-d6	14.051	166	743111	33.868	ng/u1	0.01
49) Acenaphthylene-d8	14.334	160	757277	31.700	ng/u1	0.01
54) 4-Nitrophenol-d4	14.887	143	82646	24.852	ng/u1	0.02
60) Fluorene-d10	15.640	176	649545	33.851	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.787	200	89001	22.748	ng/u1	0.02
73) Anthracene-d10	17.528	188	1078329	34.065	ng/u1	0.01
81) Pyrene-d10	19.781	212	1367641	35.831	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.980	264	1312295	34.708	ng/u1	0.02

Target Compounds Qvalue

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

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