

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011942.D
 Acq On : 05 Oct 2022 22:31
 Operator : CG/JU
 Sample : PB148103BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK103

Quant Time: Oct 05 23:25:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	240071	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	1009262	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	615786	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1301179	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1224260	20.000	ng/u1	0.00
88) Perylene-d12	23.798	264	1242960	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	37152	6.825	ng/uL	0.00
4) Pyridine-d5	3.722	84	483172	30.303	ng/u1	0.00
7) Phenol-d5	7.046	99	663619	32.020	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	379305	33.018	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	554974	33.921	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	526358	30.723	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	257769	33.531	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	275952	33.552	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	505171	33.207	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	735957	31.733	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	1507427	34.232	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1753203	34.924	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	265418	29.107	ng/u1	0.00
60) Fluorene-d10	15.522	176	1368922	35.796	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	201942	25.641	ng/u1	0.00
73) Anthracene-d10	17.386	188	2211208	37.307	ng/u1	0.00
81) Pyrene-d10	19.616	212	2536936	40.917	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	2486245	39.958	ng/u1	0.00

Target Compounds	Qvalue

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

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