

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011973.D
 Acq On : 06 Oct 2022 16:49
 Operator : CG/JU
 Sample : PB148077BL
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK077

Quant Time: Oct 06 23:07:39 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	229723	20.000	ng/ul	0.00
20) Naphthalene-d8	10.686	136	925039	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.527	164	532636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1113095	20.000	ng/ul	0.00
79) Chrysene-d12	21.368	240	1120613	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1171339	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	31931	6.130	ng/uL	0.00
4) Pyridine-d5	3.722	84	376933	24.705	ng/ul	0.00
7) Phenol-d5	7.045	99	531472	26.799	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	307422	27.966	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	446784	28.539	ng/ul	0.00
15) 4-Methylphenol-d8	8.592	113	411839	25.121	ng/ul	0.00
21) Nitrobenzene-d5	9.045	128	214007	30.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	222263	29.484	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	381371	27.351	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	555632	26.139	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	1108502	29.103	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1302477	29.996	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	186900	23.696	ng/ul	0.00
60) Fluorene-d10	15.522	176	993639	30.039	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	141751	21.040	ng/ul	0.00
73) Anthracene-d10	17.380	188	1624106	32.031	ng/ul	0.00
81) Pyrene-d10	19.615	212	1899430	33.469	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1921374	32.768	ng/ul	0.00

Target Compounds	Qvalue

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

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