

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011974.D
 Acq On : 06 Oct 2022 17:25
 Operator : CG/JU
 Sample : PB148078BL
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:07:49 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	209842	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	840584	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	474673	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	995315	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1026135	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1067001	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	31697	6.662	ng/uL	0.00
4) Pyridine-d5	3.722	84	398692	28.607	ng/ul	0.00
7) Phenol-d5	7.046	99	526818	29.081	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	311619	31.033	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	444259	31.066	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	402612	26.885	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	206528	32.257	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	214640	31.334	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	373071	29.444	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	566452	29.325	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1057127	31.143	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1250406	32.313	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	180348	25.657	ng/ul	0.00
60) Fluorene-d10	15.522	176	961540	32.618	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	134958	22.402	ng/ul	0.00
73) Anthracene-d10	17.381	188	1594602	35.171	ng/ul	0.00
81) Pyrene-d10	19.616	212	1894326	36.452	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1929131	36.117	ng/ul	0.00

Target Compounds Qvalue

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

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