

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011979.D
 Acq On : 06 Oct 2022 20:19
 Operator : CG/JU
 Sample : N4923-18
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10040

Quant Time: Oct 06 23:08:39 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	276013	20.000	ng/ul	0.00
20) Naphthalene-d8	10.686	136	1078924	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	585762	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1188213	20.000	ng/ul	0.00
79) Chrysene-d12	21.368	240	1291760	20.000	ng/ul	0.00
88) Perylene-d12	23.792	264	1338748	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	27359	4.372	ng/uL	0.00
4) Pyridine-d5	3.728	84	125783	6.862	ng/ul	0.00
7) Phenol-d5	7.045	99	117453	4.929	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	337267	25.535	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	358041	19.034	ng/ul	0.00
15) 4-Methylphenol-d8	8.592	113	183266	9.304	ng/ul	0.00
21) Nitrobenzene-d5	9.045	128	219956	26.765	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	216801	24.658	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	344023	21.154	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	340731	13.743	ng/ul	0.00
46) Dimethylphthalate-d6	13.933	166	1198417	28.610	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1361142	28.504	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	30858	3.557	ng/ul	0.02
60) Fluorene-d10	15.516	176	1078427	29.645	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	148332	20.625	ng/ul	0.00
73) Anthracene-d10	17.380	188	1674686	30.941	ng/ul	0.00
81) Pyrene-d10	19.615	212	2087324	31.906	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2209900	32.976	ng/ul	0.00

Target Compounds	Qvalue

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

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