

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011977.D
 Acq On : 06 Oct 2022 19:09
 Operator : CG/JU
 Sample : N4923-14
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10013

Quant Time: Oct 06 23:08:19 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	7.875	152	263564	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1094119	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	657366	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1427600	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1471481	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1514871	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	26389	4.416	ng/uL	0.00
4) Pyridine-d5	3.740	84	71623	4.092	ng/ul	0.02
7) Phenol-d5	7.046	99	114901	5.050	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	323793	25.673	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	353111	19.659	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	214556	11.407	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	222420	26.689	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	219708	24.641	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	377976	22.919	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	500450	19.905	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1433241	30.489	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1524619	28.450	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	36727	3.773	ng/ul	0.01
60) Fluorene-d10	15.522	176	1298929	31.818	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	201126	23.276	ng/ul	0.00
73) Anthracene-d10	17.381	188	2294040	35.277	ng/ul	0.00
81) Pyrene-d10	19.616	212	2746591	36.856	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	2762527	36.429	ng/ul	-0.01

Target Compounds Qvalue

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

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