

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011933.D
 Acq On : 05 Oct 2022 16:47
 Operator : CG/JU
 Sample : N4890-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY0E6

Quant Time: Oct 05 23:23:15 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	257248	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1021648	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	575473	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.287	188	1202781	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1226888	20.000	ng/u1	0.00
88) Perylene-d12	23.804	264	1259372	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	27624	4.736	ng/uL	0.00
4) Pyridine-d5	3.728	84	91994	5.384	ng/u1	0.00
7) Phenol-d5	7.046	99	124226	5.594	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	365588	29.699	ng/u1	0.00
11) 2-Chlorophenol-d4	7.411	132	402473	22.957	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	233999	12.746	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	243044	31.232	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	240462	28.882	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	403886	26.227	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	510666	21.752	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	1345465	32.695	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1531959	32.655	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	35774	4.198	ng/u1	0.01
60) Fluorene-d10	15.522	176	1268228	35.486	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	188163	25.846	ng/u1	0.00
73) Anthracene-d10	17.387	188	2160328	39.430	ng/u1	0.00
81) Pyrene-d10	19.622	212	2580815	41.536	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	2583115	40.974	ng/u1	0.00

Target Compounds Qvalue

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

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