

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011882.D
 Acq On : 03 Oct 2022 20:48
 Operator : CG/JU
 Sample : N4847-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8Q9

Quant Time: Oct 03 23:33:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.887	152	238728	20.000 ng/u1	0.00
20) Naphthalene-d8	10.698	136	929879	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.533	164	519482	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.292	188	1051603	20.000 ng/u1	0.00
79) Chrysene-d12	21.374	240	1207391	20.000 ng/u1	-0.01
88) Perylene-d12	23.809	264	1273909	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.305	96	26807	4.952 ng/uL	0.00
4) Pyridine-d5	3.728	84	150514	9.493 ng/u1	0.00
7) Phenol-d5	7.051	99	117579	5.705 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	301417	26.385 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	349873	21.505 ng/u1	0.00
15) 4-Methylphenol-d8	8.598	113	204477	12.002 ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	211181	29.816 ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	211306	27.885 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	346655	24.732 ng/u1	0.00
31) 4-Chloroaniline-d4	10.839	131	441780	20.675 ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1115301	30.023 ng/u1	0.00
49) Acenaphthylene-d8	14.227	160	1284336	30.327 ng/u1	0.00
54) 4-Nitrophenol-d4	14.751	143	32507	4.226 ng/u1	0.01
60) Fluorene-d10	15.527	176	995817	30.867 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	145962	22.932 ng/u1	0.00
73) Anthracene-d10	17.392	188	1643970	34.319 ng/u1	0.00
81) Pyrene-d10	19.621	212	2057900	33.655 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.650	264	2288197	35.882 ng/u1	-0.01

Target Compounds	Qvalue

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

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