

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012510.D
 Acq On : 04 Nov 2022 03:35
 Operator : CG/JU
 Sample : N5319-16
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAB1

Quant Time: Nov 04 04:35:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	332957	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1623958	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	1146717	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2651833	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	2393367	20.000	ng/u1	0.00
88) Perylene-d12	23.669	264	2144119	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	33043	3.996	ng/uL	0.00
4) Pyridine-d5	3.658	84	153749	6.473	ng/u1	0.00
7) Phenol-d5	6.958	99	209168	7.004	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	536013	30.066	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	552382	25.329	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	411063	17.404	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	370601	35.191	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	404137	36.973	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	707685	29.875	ng/u1	0.00
31) 4-Chloroaniline-d4	10.728	131	830590	23.549	ng/u1	0.00
46) Dimethylphthalate-d6	13.846	166	2528396	32.209	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2805693	31.276	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	70912	4.773	ng/u1	0.02
60) Fluorene-d10	15.428	176	2172503	32.325	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	348740	25.259	ng/u1	0.00
73) Anthracene-d10	17.292	188	3570766	31.251	ng/u1	0.00
81) Pyrene-d10	19.533	212	4090648	34.042	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.516	264	3377606	32.240	ng/u1	0.00

Target Compounds	Qvalue

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

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