

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012501.D
 Acq On : 03 Nov 2022 21:51
 Operator : CG/JU
 Sample : N5319-02
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA97

Quant Time: Nov 03 23:06:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.781	152	262246	20.000 ng/u1	0.00
20) Naphthalene-d8	10.581	136	1250614	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.434	164	860608	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.192	188	1938596	20.000 ng/u1	-0.01
79) Chrysene-d12	21.286	240	1769472	20.000 ng/u1	-0.01
88) Perylene-d12	23.669	264	1533398	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.229	96	25016	3.841 ng/uL	0.00
4) Pyridine-d5	3.664	84	134658	7.197 ng/u1	0.00
7) Phenol-d5	6.964	99	145507	6.186 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	419517	29.877 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.311	132	422263	24.583 ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	292107	15.702 ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	292556	36.074 ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	310741	36.916 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.211	165	552852	30.306 ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	666510	24.538 ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2080594	35.316 ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2180682	32.390 ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	47285	4.241 ng/u1	0.01
60) Fluorene-d10	15.428	176	1793033	35.548 ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	298204	29.545 ng/u1	0.00
73) Anthracene-d10	17.292	188	3038259	36.373 ng/u1	-0.01
81) Pyrene-d10	19.534	212	3516237	39.580 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.516	264	2800483	37.378 ng/u1	-0.01

Target Compounds Qvalue

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

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