

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012455.D
 Acq On : 02 Nov 2022 15:56
 Operator : CG/JU
 Sample : N5321-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBWR5

Quant Time: Nov 02 23:14:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.787	152	326356	20.000 ng/u1	0.00
20) Naphthalene-d8	10.587	136	1527089	20.000 ng/u1	-0.01
38) Acenaphthene-d10	14.439	164	1020800	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2659940	20.000 ng/u1	-0.01
79) Chrysene-d12	21.298	240	2447472	20.000 ng/u1	0.00
88) Perylene-d12	23.686	264	2202329	20.000 ng/u1	-0.01

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.234	96	33606	4.146 ng/uL	0.00
4) Pyridine-d5	3.664	84	218997	9.406 ng/u1	0.00
7) Phenol-d5	6.963	99	232537	7.944 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	551050	31.535 ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	580109	27.138 ng/u1	-0.01
15) 4-Methylphenol-d8	8.504	113	449253	19.405 ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	376529	38.022 ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	408897	39.782 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	741885	33.305 ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	869096	26.204 ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2755906	39.438 ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2925704	36.637 ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	85306	6.450 ng/u1	0.00
60) Fluorene-d10	15.439	176	2718426	45.437 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	480063	34.665 ng/u1	-0.01
73) Anthracene-d10	17.304	188	4475035	39.045 ng/u1	0.00
81) Pyrene-d10	19.545	212	5218393	42.468 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	4249276	39.488 ng/u1	-0.01

Target Compounds Qvalue

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

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