

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012479.D
 Acq On : 03 Nov 2022 08:01
 Operator : CG/JU
 Sample : N5321-20
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBWU9

Quant Time: Nov 03 08:29:39 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	285615	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1331801	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	908538	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2037487	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	1970352	20.000	ng/u1	0.00
88) Perylene-d12	23.674	264	1782335	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	27899	3.933	ng/uL	0.00
4) Pyridine-d5	3.664	84	111303	5.462	ng/u1	0.00
7) Phenol-d5	6.957	99	198844	7.762	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	513982	33.609	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	539682	28.848	ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	385623	19.033	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	364683	42.226	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	395494	44.120	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	699546	36.010	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	693489	23.975	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2477376	39.833	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2697630	37.955	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	69722	5.923	ng/u1	0.01
60) Fluorene-d10	15.434	176	2120709	39.826	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	379513	35.776	ng/u1	0.00
73) Anthracene-d10	17.298	188	3448597	39.282	ng/u1	0.00
81) Pyrene-d10	19.539	212	3991965	40.354	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.521	264	3475188	39.905	ng/u1	0.00

Target Compounds Qvalue

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

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