

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012499.D
 Acq On : 03 Nov 2022 20:42
 Operator : CG/JU
 Sample : N5319-01
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 03 23:06:31 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	249924	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1178738	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.433	164	818854	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	1856285	20.000	ng/ul	0.00
79) Chrysene-d12	21.292	240	1765580	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1550389	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	21349	3.440	ng/uL	0.00
4) Pyridine-d5	3.663	84	105420	5.912	ng/ul	0.00
7) Phenol-d5	6.963	99	133966	5.976	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	356118	26.612	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	368876	22.534	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	263329	14.853	ng/ul	0.00
21) Nitrobenzene-d5	8.951	128	251132	32.854	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	267285	33.689	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	463624	26.964	ng/ul	0.00
31) 4-Chloroaniline-d4	10.733	131	570595	22.288	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	1760471	31.406	ng/ul	0.00
49) Acenaphthylene-d8	14.127	160	1889739	29.500	ng/ul	0.00
54) 4-Nitrophenol-d4	14.680	143	44499	4.194	ng/ul	0.02
60) Fluorene-d10	15.433	176	1535624	31.997	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	265111	27.431	ng/ul	0.00
73) Anthracene-d10	17.292	188	2765809	34.580	ng/ul	-0.01
81) Pyrene-d10	19.533	212	3206404	36.172	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.521	264	2634018	34.771	ng/ul	0.00

Target Compounds Qvalue

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

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