

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012500.D
 Acq On : 03 Nov 2022 21:17
 Operator : CG/JU
 Sample : N5319-13
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 03 23:06:41 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	238377	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1119095	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.434	164	777406	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	1729992	20.000	ng/ul	-0.01
79) Chrysene-d12	21.292	240	1597696	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1370532	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23617	3.989	ng/uL	0.00
4) Pyridine-d5	3.664	84	137789	8.102	ng/ul	0.00
7) Phenol-d5	6.963	99	137222	6.418	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	412808	32.343	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	399202	25.568	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	279982	16.557	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	291394	40.153	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	305259	40.526	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	537911	32.952	ng/ul	0.00
31) 4-Chloroaniline-d4	10.734	131	675829	27.805	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	2047990	38.483	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	2219086	36.488	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	36155	3.590	ng/ul	0.02
60) Fluorene-d10	15.434	176	1776676	38.993	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	274707	30.499	ng/ul	0.00
73) Anthracene-d10	17.292	188	2922368	39.205	ng/ul	-0.01
81) Pyrene-d10	19.539	212	3375434	42.080	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	2647492	39.535	ng/ul	0.00

Target Compounds Qvalue

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

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