

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012355.D
 Acq On : 27 Oct 2022 22:23
 Operator : CG/JU
 Sample : N5103-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 27 23:24:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	311262	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.610	136	1444079	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	966636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	2155235	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	2265072	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	2124726	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	22018	2.848	ng/uL	0.00
4) Pyridine-d5	3.681	84	108350	4.879	ng/ul	0.00
7) Phenol-d5	6.981	99	143501	5.140	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	458093	27.487	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.340	132	440501	21.607	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	287496	13.020	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	309233	33.022	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	319721	32.894	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	542533	25.756	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	524318	16.717	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	2076794	31.385	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	2330703	30.821	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	51790	4.135	ng/ul	0.00
60) Fluorene-d10	15.457	176	1877008	33.131	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.587	200	292165	26.037	ng/ul	0.00
73) Anthracene-d10	17.316	188	3255221	35.054	ng/ul	0.00
81) Pyrene-d10	19.557	212	3982310	35.018	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3813682	36.735	ng/ul	0.00

Target Compounds	Qvalue

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

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