

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012118.D
 Acq On : 13 Oct 2022 13:35
 Operator : CG/JU
 Sample : N5013-07
 Misc :
 ALS Vial : 89 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:33:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	259203	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1038650	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.499	164	560756	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1147383	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1254724	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1354608	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	28018	4.262	ng/uL	0.00
4) Pyridine-d5	3.705	84	169662	9.704	ng/ul	0.00
7) Phenol-d5	7.017	99	169576	8.134	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	400544	33.850	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	467408	28.246	ng/ul	0.00
15) 4-Methylphenol-d8	8.564	113	293449	18.327	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	281277	38.345	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	277737	37.340	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	462388	32.076	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	611488	28.786	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1324305	37.290	ng/ul	0.00
49) Acenaphthylene-d8	14.193	160	1608937	36.637	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	40135	5.757	ng/ul	0.00
60) Fluorene-d10	15.493	176	1209057	37.629	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	168804	27.862	ng/ul	0.00
73) Anthracene-d10	17.351	188	1915514	38.446	ng/ul	0.00
81) Pyrene-d10	19.592	212	2376329	36.968	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.604	264	2566554	38.975	ng/ul	0.00

Target Compounds Qvalue

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

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