

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012110.D
 Acq On : 13 Oct 2022 08:59
 Operator : CG/JU
 Sample : N5026-09
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:30:25 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	252203	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1093354	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	693451	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1584093	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1604320	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1670674	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	25771	4.029	ng/uL	0.00
4) Pyridine-d5	3.717	84	34199	2.010	ng/ul	0.02
7) Phenol-d5	7.016	99	129965	6.407	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	315253	27.382	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	368695	22.899	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	227985	14.633	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	226254	29.301	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	230585	29.450	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	389532	25.670	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	385416	17.236	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1371956	31.239	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1532178	28.213	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	47640	5.525	ng/ul	0.00
60) Fluorene-d10	15.492	176	1279109	32.192	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	213145	25.482	ng/ul	0.00
73) Anthracene-d10	17.357	188	2296170	33.381	ng/ul	0.00
81) Pyrene-d10	19.592	212	2762998	33.617	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2732657	33.647	ng/ul	0.00

Target Compounds Qvalue

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

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