

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012123.D
 Acq On : 13 Oct 2022 16:29
 Operator : CG/JU
 Sample : N5013-02
 Misc :
 ALS Vial : 94 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:34:15 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	240275	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	967062	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	516734	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1041752	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1114369	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	1196417	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	27973	4.590	ng/uL	0.00
4) Pyridine-d5	3.699	84	179659	11.085	ng/ul	0.00
7) Phenol-d5	7.016	99	168818	8.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	396281	36.128	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	458247	29.874	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	282497	19.032	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	278379	40.759	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	276212	39.884	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	449604	33.498	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	641162	32.417	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1286372	39.308	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1570016	38.797	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	37071	5.770	ng/ul	0.00
60) Fluorene-d10	15.492	176	1160713	39.202	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	156332	28.420	ng/ul	0.00
73) Anthracene-d10	17.351	188	1803531	39.869	ng/ul	0.00
81) Pyrene-d10	19.592	212	2195980	38.465	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2367311	40.703	ng/ul	0.00

Target Compounds Qvalue

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

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