

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012120.D
 Acq On : 13 Oct 2022 14:44
 Operator : CG/JU
 Sample : N5041-02
 Misc :
 ALS Vial : 91 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:33:39 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	7.852	152	271936	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1097558	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	602013	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1223969	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1314114	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	1388320	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	28427	4.121	ng/uL	0.00
4) Pyridine-d5	3.711	84	49604	2.704	ng/ul	0.01
7) Phenol-d5	7.016	99	166705	7.622	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	383702	30.908	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	452165	26.046	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	282303	16.805	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	267623	34.525	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	274490	34.923	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	442175	29.028	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	343488	15.302	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1311306	34.394	ng/ul	0.00
49) Acenaphthylene-d8	14.192	160	1566408	33.224	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	40577	5.421	ng/ul	0.00
60) Fluorene-d10	15.492	176	1202280	34.854	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	175369	27.135	ng/ul	0.00
73) Anthracene-d10	17.357	188	1995682	37.549	ng/ul	0.00
81) Pyrene-d10	19.592	212	2398533	35.627	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2569236	38.069	ng/ul	0.00

Target Compounds Qvalue

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

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