

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012135.D
 Acq On : 14 Oct 2022 13:32
 Operator : CG/JU
 Sample : N4982-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 14 23:03:02 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	191722	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	796627	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	421254	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	801883	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	867059	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	931567	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	24462	5.030	ng/uL	0.00
4) Pyridine-d5	3.705	84	87595	6.773	ng/ul	0.00
7) Phenol-d5	7.016	99	137134	8.894	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	357641	40.863	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	393709	32.167	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	215301	18.179	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	241248	42.879	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	238041	41.726	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	370738	33.532	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	320544	19.674	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1074681	40.282	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	1310944	39.737	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	23175	4.425	ng/ul	0.01
60) Fluorene-d10	15.492	176	972978	40.310	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	115391	27.252	ng/ul	0.00
73) Anthracene-d10	17.351	188	1451482	41.684	ng/ul	0.00
81) Pyrene-d10	19.592	212	1807303	40.687	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.609	264	1988736	43.915	ng/ul	0.00

Target Compounds Qvalue

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

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