

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012146.D
 Acq On : 14 Oct 2022 20:27
 Operator : CG/JU
 Sample : PB148313BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 14 23:06:10 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	7.840	152	271742	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.646	136	1099007	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.492	164	592078	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1179655	20.000	ng/ul	0.00
79) Chrysene-d12	21.339	240	1234703	20.000	ng/ul	# 0.00
88) Perylene-d12	23.757	264	1308452	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	24957	3.621	ng/uL	0.00
4) Pyridine-d5	3.693	84	348640	19.021	ng/ul	0.00
7) Phenol-d5	7.010	99	598980	27.407	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	352601	28.423	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.375	132	481340	27.746	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113	448862	26.739	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	227394	29.297	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	232914	29.594	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.269	165	395048	25.900	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.793	131	588362	26.176	ng/ul	-0.01
46) Dimethylphthalate-d6	13.904	166	1059441	28.254	ng/ul	-0.01
49) Acenaphthylene-d8	14.187	160	1303462	28.111	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	164388	22.331	ng/ul	0.00
60) Fluorene-d10	15.487	176	964631	28.434	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.616	200	115228	18.499	ng/ul	0.00
73) Anthracene-d10	17.351	188	1545079	30.163	ng/ul	0.00
81) Pyrene-d10	19.586	212	1812938	28.661	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.598	264	1907014	29.981	ng/ul	-0.01

Target Compounds Qvalue

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

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