

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012065.D
 Acq On : 12 Oct 2022 06:00
 Operator : CG/JU
 Sample : N4991-08
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBG0

Quant Time: Oct 12 07:25:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	234944	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	959138	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.498	164	557220	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.257	188	1214854	20.000	ng/u1	-0.01
79) Chrysene-d12	21.345	240	1315343	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	1401482	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	32951	5.530	ng/uL	0.00
4) Pyridine-d5	3.699	84	358036	22.593	ng/u1	0.00
7) Phenol-d5	7.022	99	445764	23.591	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	350511	32.680	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	449017	29.937	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	350604	24.157	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	252792	37.318	ng/u1	0.00
24) 2-Nitrophenol-d4	9.746	143	270543	39.388	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	461894	34.698	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	579508	29.542	ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	1535499	43.511	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	1849222	42.376	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	159207	22.980	ng/u1	0.00
60) Fluorene-d10	15.498	176	1425766	44.655	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	183770	28.648	ng/u1	0.00
73) Anthracene-d10	17.357	188	2268426	43.000	ng/u1	0.00
81) Pyrene-d10	19.598	212	2802486	41.589	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	2877532	42.236	ng/u1	0.00

Target Compounds Qvalue

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

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