

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012422.D
 Acq On : 01 Nov 2022 12:07
 Operator : CG/JU
 Sample : N5270-06
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA81

Quant Time: Nov 01 22:56:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	336973	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1523786	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.440	164	995264	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2182839	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	2343807	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	2313445	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	41502	4.959	ng/uL	0.00
4) Pyridine-d5	3.664	84	249733	10.388	ng/u1	0.00
7) Phenol-d5	6.964	99	239061	7.910	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	657338	36.432	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	652614	29.568	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	460299	19.256	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	444776	45.011	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	463168	45.160	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	773284	34.790	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	1062677	32.110	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2749319	40.353	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	3128582	40.183	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	84061	6.519	ng/u1	0.00
60) Fluorene-d10	15.440	176	2407128	41.266	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	420873	37.033	ng/u1	0.00
73) Anthracene-d10	17.304	188	3880651	41.260	ng/u1	0.00
81) Pyrene-d10	19.545	212	4645829	39.480	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	4692715	41.514	ng/u1	-0.01

Target Compounds	Qvalue

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

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