

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011518.D
 Acq On : 19 Aug 2022 21:50
 Operator : CG/JU
 Sample : N4213-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Aug 20 04:41:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	88209	20.000	ng/u1	0.00
20) Naphthalene-d8	10.340	136	369873	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.228	164	203153	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.998	188	403109	20.000	ng/u1	0.00
79) Chrysene-d12	21.122	240	388573	20.000	ng/u1	-0.01
88) Perylene-d12	23.410	264	412213	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	16813	5.915	ng/uL	0.00
4) Pyridine-d5	3.481	84	64703	8.891	ng/u1	0.00
7) Phenol-d5	6.740	99	257589	31.777	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	172387	32.734	ng/u1	0.00
11) 2-Chlorophenol-d4	7.087	132	210626	35.328	ng/u1	0.00
15) 4-Methylphenol-d8	8.275	113	205729	31.709	ng/u1	0.00
21) Nitrobenzene-d5	8.716	128	101081	37.960	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.434	143	103422	38.163	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	172716	34.527	ng/u1	0.00
31) 4-Chloroaniline-d4	10.493	131	249497	29.962	ng/u1	0.00
46) Dimethylphthalate-d6	13.651	166	531009	36.956	ng/u1	0.00
49) Acenaphthylene-d8	13.916	160	606841	37.999	ng/u1	0.00
54) 4-Nitrophenol-d4	14.457	143	79957	28.882	ng/u1	0.00
60) Fluorene-d10	15.228	176	447901	37.053	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	67433	30.958	ng/u1	0.00
73) Anthracene-d10	17.098	188	685233	38.382	ng/u1	0.00
81) Pyrene-d10	19.357	212	759547	38.635	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.269	264	780151	38.907	ng/u1	-0.01

Target Compounds Qvalue

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

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