

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012419.D
 Acq On : 01 Nov 2022 10:24
 Operator : CG/JU
 Sample : N5270-12
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 01 22:56:24 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	423862	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1889125	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.445	164	1234701	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2708753	20.000	ng/u1	0.00
79) Chrysene-d12	21.304	240	2842505	20.000	ng/u1	0.00
88) Perylene-d12	23.692	264	2724308	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	45829	4.354	ng/uL	0.00
4) Pyridine-d5	3.670	84	115749	3.828	ng/u1	0.01
7) Phenol-d5	6.964	99	298716	7.857	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	749205	33.012	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	762777	27.475	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	557894	18.554	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	512014	41.795	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	538782	42.373	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	911104	33.063	ng/u1	0.00
31) 4-Chloroaniline-d4	10.746	131	584760	14.252	ng/u1	0.00
46) Dimethylphthalate-d6	13.863	166	3506410	41.485	ng/u1	0.00
49) Acenaphthylene-d8	14.140	160	3732529	38.643	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	139654	8.730	ng/u1	0.00
60) Fluorene-d10	15.440	176	3020680	41.742	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	586238	41.569	ng/u1	0.00
73) Anthracene-d10	17.304	188	4997416	42.818	ng/u1	0.00
81) Pyrene-d10	19.551	212	5899161	41.336	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	5791085	43.505	ng/u1	0.00

Target Compounds	Qvalue

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

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