

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012433.D
 Acq On : 01 Nov 2022 18:27
 Operator : CG/JU
 Sample : N5270-01
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 01 23:00:46 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	339836	20.000	ng/ul	0.00
20) Naphthalene-d8	10.592	136	1551242	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	1026592	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	2239216	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	2177322	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1912530	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	36094	4.277	ng/uL	0.00
4) Pyridine-d5	3.663	84	217944	8.989	ng/ul	0.00
7) Phenol-d5	6.963	99	198429	6.510	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	584635	32.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	545978	24.529	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	376589	15.621	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	396193	39.385	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	396615	37.986	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	648633	28.666	ng/ul	0.00
31) 4-Chloroaniline-d4	10.739	131	931900	27.660	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2490032	35.432	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	2812341	35.019	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	54437	4.093	ng/ul	0.00
60) Fluorene-d10	15.439	176	2182794	36.278	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	344264	29.529	ng/ul	0.00
73) Anthracene-d10	17.304	188	3507478	36.353	ng/ul	0.00
81) Pyrene-d10	19.545	212	4081691	37.338	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	3447417	36.891	ng/ul	-0.01

Target Compounds	Qvalue

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

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