

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
 Data File : BP012943.D
 Acq On : 02 Dec 2022 17:09
 Operator : CG/JU
 Sample : N5785-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COL73

Quant Time: Dec 03 01:31:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 01 23:46:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	300153	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	1543107	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	1253796	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	3076214	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	3511154	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	3455260	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	22454	3.329	ng/uL	0.00
4) Pyridine-d5	3.828	84	122272	5.617	ng/ul	0.01
7) Phenol-d5	7.158	99	113023	4.078	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	384083	24.475	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	351388	17.296	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	253597	11.142	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	267759	24.400	ng/ul	0.00
24) 2-Nitrophenol-d4	9.899	143	126586	10.080	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	472473	18.793	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	390696	10.999	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	2261481	25.008	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	2498609	25.025	ng/ul	0.00
54) 4-Nitrophenol-d4	14.839	143	3443	0.186	ng/ul	0.01
60) Fluorene-d10	15.634	176	2113977	26.596	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	11505	0.637	ng/ul	0.00
73) Anthracene-d10	17.498	188	4288547	31.540	ng/ul	0.00
81) Pyrene-d10	19.722	212	5263153	30.495	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	5203382	30.821	ng/ul	0.00

Target Compounds Qvalue

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

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