

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012231.D
 Acq On : 21 Oct 2022 01:05
 Operator : CG/JU
 Sample : N5103-09
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBB1

Quant Time: Oct 21 01:43:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	416018	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1834949	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1170676	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2387069	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1815694	20.000	ng/u1	0.00
88) Perylene-d12	23.727	264	1734129	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	37775	3.675	ng/uL	0.00
4) Pyridine-d5	3.687	84	182728	6.169	ng/u1	0.00
7) Phenol-d5	6.993	99	198803	5.535	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	619304	28.433	ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	610340	22.326	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	391957	14.042	ng/u1	-0.01
21) Nitrobenzene-d5	8.993	128	420193	30.936	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	444280	30.317	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	738422	26.864	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	606087	15.519	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	2651609	32.562	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2969961	31.951	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	70732	4.636	ng/u1	0.00
60) Fluorene-d10	15.469	176	2314209	33.033	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	358329	26.387	ng/u1	0.00
73) Anthracene-d10	17.333	188	3667996	35.056	ng/u1	0.00
81) Pyrene-d10	19.569	212	3911938	37.841	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.574	264	3166158	36.408	ng/u1	-0.01

Target Compounds Qvalue

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

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