

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012227.D
 Acq On : 20 Oct 2022 22:49
 Operator : CG/JU
 Sample : N5103-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBX2

Quant Time: Oct 21 01:26:46 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	376935	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1663358	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1046517	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2096870	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1606259	20.000	ng/u1	0.00
88) Perylene-d12	23.727	264	1485441	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	44024	4.727	ng/uL	0.00
4) Pyridine-d5	3.687	84	209805	7.817	ng/u1	0.00
7) Phenol-d5	6.993	99	205436	6.313	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	637923	32.324	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	617022	24.910	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	385451	15.240	ng/u1	0.00
21) Nitrobenzene-d5	8.992	128	438522	35.617	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	452858	34.091	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	734483	29.477	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	795852	22.480	ng/u1	0.00
46) Dimethylphthalate-d6	13.886	166	2732621	37.537	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2919366	35.133	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	74373	5.453	ng/u1	0.00
60) Fluorene-d10	15.469	176	2338194	37.335	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	353682	29.649	ng/u1	0.00
73) Anthracene-d10	17.333	188	3771130	41.030	ng/u1	0.00
81) Pyrene-d10	19.574	212	4035046	44.122	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	3151902	42.312	ng/u1	-0.01

Target Compounds Qvalue

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

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