

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012228.D
 Acq On : 20 Oct 2022 23:23
 Operator : CG/JU
 Sample : N5103-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBX0

Quant Time: Oct 21 01:28:10 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	445089	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1915678	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1166733	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2203758	20.000	ng/u1	0.00
79) Chrysene-d12	21.328	240	1562511	20.000	ng/u1	0.00
88) Perylene-d12	23.727	264	1602361	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	43099	3.919	ng/uL	0.00
4) Pyridine-d5	3.687	84	205629	6.489	ng/u1	0.00
7) Phenol-d5	6.993	99	204336	5.318	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	609058	26.136	ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	601775	20.575	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	397943	13.325	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	417200	29.422	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	433111	28.310	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	716026	24.951	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	851755	20.890	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	2691704	33.166	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	2862265	30.897	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.693	143	69782	4.589	ng/u1	0.00
60) Fluorene-d10	15.469	176	2268889	32.495	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	316454	25.242	ng/u1	0.00
73) Anthracene-d10	17.334	188	3645109	37.735	ng/u1	0.00
81) Pyrene-d10	19.575	212	3597452	40.438	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	3138299	39.055	ng/u1	-0.01

Target Compounds Qvalue

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

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