

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012215.D
 Acq On : 20 Oct 2022 15:57
 Operator : CG/JU
 Sample : N5103-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 20 23:23:30 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	378889	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1644912	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1027340	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.239	188	2032474	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1449258	20.000	ng/u1	0.00
88) Perylene-d12	23.739	264	1445422	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	37855	4.043	ng/uL	0.00
4) Pyridine-d5	3.687	84	226806	8.407	ng/u1	0.00
7) Phenol-d5	6.993	99	168909	5.164	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	599343	30.213	ng/u1	0.00
11) 2-Chlorophenol-d4	7.357	132	554883	22.286	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	336912	13.253	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	373247	30.655	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	378047	28.778	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	660732	26.814	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	451778	12.904	ng/u1	0.00
46) Dimethylphthalate-d6	13.892	166	2427298	33.966	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2725810	33.416	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	38319	2.862	ng/u1	0.00
60) Fluorene-d10	15.475	176	2155937	35.067	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	318298	27.528	ng/u1	0.00
73) Anthracene-d10	17.333	188	3216722	36.107	ng/u1	0.00
81) Pyrene-d10	19.575	212	3238868	39.252	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	2641760	36.445	ng/u1	0.00

Target Compounds Qvalue

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

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