

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042318.D
 Acq On : 17 Oct 2023 11:43
 Operator : MA/JU
 Sample : 04900-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW2

Quant Time: Oct 17 13:39:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 15:38:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	135293	20.000	ng/ul	0.00
20) Naphthalene-d8	10.827	136	616416	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	373590	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	792902	20.000	ng/ul	0.00
79) Chrysene-d12	21.586	240	710642	20.000	ng/ul	-0.01
88) Perylene-d12	24.062	264	767957	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	22025	5.087	ng/uL	0.00
4) Pyridine-d5	3.804	84	97331	7.590	ng/ul	0.00
7) Phenol-d5	7.151	99	100216	6.684	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	378023	37.313	ng/ul	0.00
11) 2-Chlorophenol-d4	7.528	132	285025	27.633	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	189451	15.955	ng/ul	0.00
21) Nitrobenzene-d5	9.204	128	191901	36.657	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	195556	34.944	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	313623	30.709	ng/ul	0.00
31) 4-Chloroaniline-d4	10.992	131	414336	25.767	ng/ul	0.00
46) Dimethylphthalate-d6	14.068	166	1201350	37.096	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1324109	37.357	ng/ul	0.00
54) 4-Nitrophenol-d4	14.862	143	30888	4.742	ng/ul	-0.01
60) Fluorene-d10	15.645	176	1022456	38.406	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.768	200	150060	27.017	ng/ul	0.00
73) Anthracene-d10	17.503	188	1608011	39.589	ng/ul	0.00
81) Pyrene-d10	19.797	212	1899185	43.849	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.903	264	1758237	41.711	ng/ul	-0.01

Target Compounds Qvalue

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

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