

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044987.D
 Acq On : 07 Apr 2024 01:38
 Operator : MA/JU
 Sample : P2013-03
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Apr 08 05:05:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 05:01:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	83594	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	382173	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	241793	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.039	188	511068	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240	499186	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	589978	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	10274	4.589	ng/uL	0.00
4) Pyridine-d5	3.622	84	34932	5.738	ng/ul	0.01
7) Phenol-d5	6.816	99	49718	7.016	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	136218	32.263	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	152244	27.421	ng/ul	0.00
15) 4-Methylphenol-d8	8.345	113	99919	18.013	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	94772	32.284	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	100319	32.377	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	177323	31.214	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	160091	17.700	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	631046	37.483	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160	706083	35.501	ng/ul	0.00
54) 4-Nitrophenol-d4	14.533	143	18511	5.268	ng/ul	0.02
60) Fluorene-d10	15.280	176	557177	37.123	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	87594	28.907	ng/ul	0.00
73) Anthracene-d10	17.139	188	913950	39.753	ng/ul	0.00
81) Pyrene-d10	19.445	212	1099809	44.744	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.333	264	1174275	40.801	ng/ul	0.00

Target Compounds Qvalue

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

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