

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044993.D
 Acq On : 07 Apr 2024 05:13
 Operator : MA/JU
 Sample : P2013-09
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0MB5

Quant Time: Apr 08 05:07:20 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	78201	20.000	ng/u1	0.00
20) Naphthalene-d8	10.404	136	347512	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.280	164	214044	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.039	188	445304	20.000	ng/u1	0.00
79) Chrysene-d12	21.244	240	437571	20.000	ng/u1	0.00
88) Perylene-d12	23.468	264	524551	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	12169	5.811	ng/uL	0.00
4) Pyridine-d5	3.616	84	64917	11.399	ng/u1	0.00
7) Phenol-d5	6.816	99	69040	10.414	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	183277	46.402	ng/u1	0.00
11) 2-Chlorophenol-d4	7.169	132	201554	38.806	ng/u1	0.00
15) 4-Methylphenol-d8	8.345	113	124688	24.028	ng/u1	0.00
21) Nitrobenzene-d5	8.792	128	123773	46.368	ng/u1	0.00
24) 2-Nitrophenol-d4	9.504	143	135405	48.059	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	226653	43.877	ng/u1	0.00
31) 4-Chloroaniline-d4	10.563	131	183303	22.288	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	712425	47.803	ng/u1	0.00
49) Acenaphthylene-d8	13.974	160	848341	48.184	ng/u1	0.00
54) 4-Nitrophenol-d4	14.533	143	19870	6.388	ng/u1	0.02
60) Fluorene-d10	15.280	176	640175	48.182	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	86923	32.922	ng/u1	0.00
73) Anthracene-d10	17.139	188	992753	49.558	ng/u1	0.00
81) Pyrene-d10	19.445	212	1201544	55.767	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.333	264	1311067	51.235	ng/u1	0.00

Target Compounds Qvalue

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

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