

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034171.D
 Acq On : 09 Mar 2022 04:55
 Operator : CG/JU
 Sample : N1733-19
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRT7

Quant Time: Mar 09 05:53:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 13:09:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.989	152	160109	20.000	ng/u1	0.00
20) Naphthalene-d8	10.801	136	597586	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.624	164	385697	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.365	188	796884	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.536	240	836371	20.000	ng/u1	# 0.00
88) Perylene-d12	23.983	264	937241	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	25926	7.448	ng/uL	0.00
4) Pyridine-d5	3.790	84	245822	24.258	ng/u1	0.00
7) Phenol-d5	7.137	99	431824	34.421	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	243816	36.108	ng/u1	0.00
11) 2-Chlorophenol-d4	7.519	132	385311	37.051	ng/u1	0.00
15) 4-Methylphenol-d8	8.689	113	355008	33.772	ng/u1	0.00
21) Nitrobenzene-d5	9.148	128	180805	39.049	ng/u1	0.00
24) 2-Nitrophenol-d4	9.878	143	194947	38.765	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	379807	36.623	ng/u1	0.00
31) 4-Chloroaniline-d4	10.930	131	476293	33.614	ng/u1	0.00
46) Dimethylphthalate-d6	14.030	166	1070434	36.573	ng/u1	0.00
49) Acenaphthylene-d8	14.318	160	1384774	37.836	ng/u1	0.00
54) 4-Nitrophenol-d4	14.801	143	158345	29.076	ng/u1	0.00
60) Fluorene-d10	15.613	176	971131	37.282	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	140287	27.280	ng/u1	0.00
73) Anthracene-d10	17.465	188	1514237	38.575	ng/u1	0.00
81) Pyrene-d10	19.753	212	1746109	37.981	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.824	264	1925110	38.720	ng/u1	0.00

Target Compounds Qvalue

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

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