

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034165.D
 Acq On : 09 Mar 2022 01:19
 Operator : CG/JU
 Sample : N1733-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRS1

Quant Time: Mar 09 02:31:54 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	7.990	152	151875	20.000	ng/u1	0.00
20) Naphthalene-d8	10.807	136	570350	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.625	164	353846	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.366	188	686898	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.536	240	681972	20.000	ng/u1	# 0.00
88) Perylene-d12	23.977	264	761737	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	21174	6.412	ng/uL	0.00
4) Pyridine-d5	3.790	84	279074	29.033	ng/u1	0.00
7) Phenol-d5	7.137	99	379190	31.864	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	212086	33.112	ng/u1	0.00
11) 2-Chlorophenol-d4	7.519	132	334824	33.942	ng/u1	0.00
15) 4-Methylphenol-d8	8.695	113	316520	31.743	ng/u1	0.00
21) Nitrobenzene-d5	9.154	128	161663	36.582	ng/u1	0.00
24) 2-Nitrophenol-d4	9.878	143	182261	37.973	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	337062	34.053	ng/u1	0.00
31) 4-Chloroaniline-d4	10.931	131	430029	31.798	ng/u1	0.00
46) Dimethylphthalate-d6	14.030	166	953179	35.499	ng/u1	0.00
49) Acenaphthylene-d8	14.319	160	1204090	35.861	ng/u1	0.00
54) 4-Nitrophenol-d4	14.801	143	133204	26.661	ng/u1	0.00
60) Fluorene-d10	15.619	176	835532	34.963	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	128340	28.952	ng/u1	0.00
73) Anthracene-d10	17.466	188	1244304	36.774	ng/u1	0.00
81) Pyrene-d10	19.754	212	1413300	37.702	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.824	264	1513757	37.462	ng/u1	0.00

Target Compounds Qvalue

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

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