

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032522\
 Data File : BM034383.D
 Acq On : 25 Mar 2022 13:59
 Operator : CG/JU
 Sample : N2084-01DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BLDG-3DL

Quant Time: Mar 26 05:38:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 02:40:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	156341	20.000	ng	0.00
21) Naphthalene-d8	10.754	136	617561	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	366609	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	700619	20.000	ng	0.00
76) Chrysene-d12	21.500	240	666780	20.000	ng	0.00
86) Perylene-d12	23.918	264	775055	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	87271	10.014	ng	0.00
7) Phenol-d6	7.101	99	107022	8.577	ng	0.00
23) Nitrobenzene-d5	9.101	82	73157	5.767	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	39899	8.074	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	156286	5.825	ng	0.00
79) Terphenyl-d14	19.942	244	224351	5.847	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.807	128	1541449	47.911	ng	100
37) 2-Methylnaphthalene	12.413	142	63127	2.750	ng	99
71) Phenanthrene	17.365	178	79987	2.039	ng	99

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

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