

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033121\
 Data File : BN014150.D
 Acq On : 31 Mar 2021 19:28
 Operator : CG/JU
 Sample : M1842-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE104D2-20210323

Quant Time: Apr 01 01:24:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN032621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 31 12:15:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.692	152	5878	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	21403	0.40	ng	# 0.00
13) Acenaphthene-d10	14.321	164	11540	0.40	ng	0.00
19) Phenanthrene-d10	17.067	188	22875	0.40	ng	0.00
29) Chrysene-d12	21.247	240	21985	0.40	ng	0.00
36) Perylene-d12	23.466	264	21351	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.292	112	1728	0.11	ng	0.00
5) Phenol-d6	6.855	99	1183	0.06	ng	0.00
8) Nitrobenzene-d5	8.832	82	4381	0.30	ng	0.00
11) 2-Methylnaphthalene-d10	12.057	152	9751	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.818	330	934	0.25	ng	0.01
15) 2-Fluorobiphenyl	12.938	172	13361	0.31	ng	0.00
27) Fluoranthene-d10	19.096	212	25164	0.42	ng	0.00
31) Terphenyl-d14	19.697	244	31879	0.63	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.231	88	8870	1.17	ng	# 87
6) bis(2-Chloroethyl)ether	7.120	93	2797	0.18	ng	96
25) Phenanthrene	17.104	178	1681	0.03	ng	# 93
34) Bis(2-ethylhexyl)phtha...	21.173	149	1930	0.10	ng	96

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

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