

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062321\
 Data File : BN014980.D
 Acq On : 22 Jun 2021 19:04
 Operator : CG/JU
 Sample : M2750-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE120D2-20210615

Manual Integrations
 APPROVED

mohammad
 6/23/2021 4:50:22 PM

Quant Time: Jun 23 03:39:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN061621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 17 01:26:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	20431	0.400	ng	#-0.01
7) Naphthalene-d8	10.518	136	92241	0.400	ng	0.00
13) Acenaphthene-d10	14.358	164	48895	0.400	ng	-0.01
19) Phenanthrene-d10	17.103	188	94140	0.400	ng	-0.01
29) Chrysene-d12	21.302	240	82565	0.400	ng	#-0.02
35) Perylene-d12	23.560	264	71162	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.380	112	9032	0.143	ng	0.00
5) Phenol-d6	6.929	99	6212	0.078	ng	0.00
8) Nitrobenzene-d5	8.883	82	24199	0.347	ng	-0.02
11) 2-Methylnaphthalene-d10	12.106	152	53682	0.339	ng	-0.01
14) 2,4,6-Tribromophenol	15.855	330	6853	0.299	ng	-0.01
15) 2-Fluorobiphenyl	12.988	172	81910	0.441	ng	-0.01
27) Fluoranthene-d10	19.143	212	106978	0.360	ng	-0.01
31) Terphenyl-d14	19.754	244	128351	0.627	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.361	88	69309m	8.089	ng	
6) bis(2-Chloroethyl)ether	7.187	93	1446	0.023	ng	# 86
34) Bis(2-ethylhexyl)phtha...	21.239	149	25830	0.119	ng	# 87

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

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