

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062421\
 Data File : BN015034.D
 Acq On : 24 Jun 2021 07:58
 Operator : CG/JU
 Sample : M2784-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE137-20210618

Manual Integrations
 APPROVED

mohammad
 6/24/2021 9:53:43 AM

Quant Time: Jun 24 08:38:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN061621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 23 16:24:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	22205	0.400	ng	# 0.00
7) Naphthalene-d8	10.505	136	99764	0.400	ng	0.00
13) Acenaphthene-d10	14.358	164	52720	0.400	ng	0.00
19) Phenanthrene-d10	17.104	188	102217	0.400	ng	# 0.00
29) Chrysene-d12	21.292	240	87932	0.400	ng	#-0.01
35) Perylene-d12	23.543	264	74483	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.380	112	11996	0.175	ng	0.00
5) Phenol-d6	6.920	99	8526	0.098	ng	0.00
8) Nitrobenzene-d5	8.870	82	28579	0.379	ng	-0.01
11) 2-Methylnaphthalene-d10	12.097	152	60097	0.351	ng	0.00
14) 2,4,6-Tribromophenol	15.843	330	7494	0.303	ng	0.00
15) 2-Fluorobiphenyl	12.975	172	88408	0.441	ng	0.00
27) Fluoranthene-d10	19.133	212	116620	0.362	ng	0.00
31) Terphenyl-d14	19.739	244	103462	0.475	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.361	88	64054m	6.878	ng	
6) bis(2-Chloroethyl)ether	7.178	93	3738	0.055	ng	# 84
34) Bis(2-ethylhexyl)phtha...	21.228	149	59383	0.257	ng	# 87

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

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