

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070121\
 Data File : BN015275.D
 Acq On : 02 Jul 2021 00:08
 Operator : CG/JU
 Sample : M2783-21DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE132D5-20210618DL

Manual Integrations
 APPROVED

mohammad
 7/2/2021 10:39:05 AM

Quant Time: Jul 02 01:01:11 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 09:06:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.734	152	25326	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	117557	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	57472	0.40	ng	0.00
19) Phenanthrene-d10	17.090	188	124238	0.40	ng	0.00
29) Chrysene-d12	21.291	240	95895	0.40	ng	#-0.01
35) Perylene-d12	23.539	264	34181	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	4420	0.07	ng	0.00
5) Phenol-d6	6.905	99	2909	0.04	ng	0.00
8) Nitrobenzene-d5	8.873	82	13455	0.20	ng	0.00
11) 2-Methylnaphthalene-d10	12.087	152	30035	0.15	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	3335	0.35	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	44533	0.19	ng	0.00
27) Fluoranthene-d10	19.127	212	59248	0.16	ng	0.00
31) Terphenyl-d14	19.733	244	53992	0.24	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	30614m	2.58	ng	
6) bis(2-Chloroethyl)ether	7.163	93	1946	0.02	ng	# 73
24) Pentachlorophenol	16.750	266	57	0.25	ng	98
34) Bis(2-ethylhexyl)phtha...	21.217	149	21253	0.33	ng	# 88

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

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