

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092521\
 Data File : BN016478.D
 Acq On : 25 Sep 2021 16:56
 Operator : CG/JU
 Sample : M3782-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE132D7-20210914

Manual Integrations
 APPROVED

mohammad
 9/28/2021 1:36:45 PM

Quant Time: Sep 26 04:15:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 03:35:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	22035	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	94747	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	48076	0.40	ng	0.00
19) Phenanthrene-d10	17.042	188	96386	0.40	ng	0.00
29) Chrysene-d12	21.249	240	88868	0.40	ng	# 0.00
35) Perylene-d12	23.519	264	77915	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.291	112	5621	0.11	ng	0.00
5) Phenol-d6	6.849	99	3705	0.06	ng	0.00
8) Nitrobenzene-d5	8.810	82	18071	0.27	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	36791	0.26	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	3521	0.32	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	49234	0.24	ng	-0.01
27) Fluoranthene-d10	19.088	212	101461	0.38	ng	0.00
31) Terphenyl-d14	19.698	244	85747	0.42	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.281	88	16108m	1.71	ng	
6) bis(2-Chloroethyl)ether	7.108	93	1715	0.03	ng	92
34) Bis(2-ethylhexyl)phtha...	21.185	149	17330	0.25	ng	97

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

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