

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092421\
 Data File : BN016461.D
 Acq On : 25 Sep 2021 06:37
 Operator : CG/JU
 Sample : M3782-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW181S-20210914

Manual Integrations
 APPROVED

mohammad
 9/27/2021 3:03:25 PM

Quant Time: Sep 25 07:37:23 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	21401	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	94645	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	49238	0.40	ng	0.00
19) Phenanthrene-d10	17.042	188	99696	0.40	ng	0.00
29) Chrysene-d12	21.249	240	95048	0.40	ng	# 0.00
35) Perylene-d12	23.515	264	82733	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	6765	0.14	ng	0.02
5) Phenol-d6	6.849	99	4627	0.08	ng	0.00
8) Nitrobenzene-d5	8.810	82	21424	0.32	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	45185	0.32	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	4691	0.38	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	60178	0.29	ng	-0.01
27) Fluoranthene-d10	19.088	212	111836	0.41	ng	0.00
31) Terphenyl-d14	19.698	244	94977	0.43	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.290	88	1658m	0.18	ng	
24) Pentachlorophenol	16.701	266	58	0.25	ng	87
34) Bis(2-ethylhexyl)phtha...	21.185	149	25649	0.34	ng	97

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

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