

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092421\
 Data File : BN016437.D
 Acq On : 24 Sep 2021 12:58
 Operator : CG/JU
 Sample : M3782-20
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP05-20210915

Manual Integrations
 APPROVED

mohammad
 9/27/2021 2:59:55 PM

Quant Time: Sep 24 13:46:21 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	22800	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	102602	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	52370	0.40	ng	0.00
19) Phenanthrene-d10	17.042	188	106598	0.40	ng	0.00
29) Chrysene-d12	21.259	240	100344	0.40	ng	# 0.01
35) Perylene-d12	23.518	264	75932	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	6012	0.12	ng	0.02
5) Phenol-d6	6.849	99	3912	0.06	ng	0.00
8) Nitrobenzene-d5	8.809	82	16605	0.23	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	46415	0.31	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	3582	0.31	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	67070	0.30	ng	-0.01
27) Fluoranthene-d10	19.087	212	117977	0.40	ng	0.00
31) Terphenyl-d14	19.698	244	127220	0.55	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.290	88	39094m	4.00	ng	
6) bis(2-Chloroethyl)ether	7.107	93	2234	0.04	ng	91
24) Pentachlorophenol	16.701	266	280	0.25	ng	92
34) Bis(2-ethylhexyl)phtha...	21.195	149	12081	0.15	ng	98

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

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