

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092021\
 Data File : BN016364.D
 Acq On : 21 Sep 2021 04:22
 Operator : CG/JU
 Sample : M3781-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW180D-20210915

Manual Integrations
 APPROVED

mohammad
 9/21/2021 11:30:48 AM

Quant Time: Sep 21 05:00:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 20 11:09:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.679	152	24838	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	114076	0.400	ng	#-0.01
13) Acenaphthene-d10	14.307	164	61587	0.400	ng	0.00
19) Phenanthrene-d10	17.054	188	129028	0.400	ng	0.00
29) Chrysene-d12	21.249	240	134404	0.400	ng	#-0.01
35) Perylene-d12	23.515	264	122698	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.310	112	7783	0.125	ng	0.00
5) Phenol-d6	6.868	99	5257	0.070	ng	0.00
8) Nitrobenzene-d5	8.822	82	24784	0.366	ng	0.00
11) 2-Methylnaphthalene-d10	12.043	152	53599	0.298	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	6045	0.334	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	74221	0.308	ng	-0.01
27) Fluoranthene-d10	19.093	212	146342	0.392	ng	0.00
31) Terphenyl-d14	19.703	244	134401	0.432	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.299	88	1465	0.142	ng	# 81
25) Phenanthrene	17.091	178	15546m	0.043	ng	
34) Bis(2-ethylhexyl)phtha...	21.196	149	64666	0.278	ng	98

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

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