

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092221\
 Data File : BN016408.D
 Acq On : 22 Sep 2021 12:23
 Operator : CG/JU
 Sample : M3781-11DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE134D2-20210916DL

Quant Time: Sep 22 12:54:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 13:21:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	24605	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	108036	0.400	ng	0.00
13) Acenaphthene-d10	14.294	164	57481	0.400	ng	-0.01
19) Phenanthrene-d10	17.054	188	112638	0.400	ng	# 0.00
29) Chrysene-d12	21.259	240	99907	0.400	ng	# 0.00
35) Perylene-d12	23.519	264	92218	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	553	0.009	ng	0.02
5) Phenol-d6	6.859	99	458	0.006	ng	0.00
8) Nitrobenzene-d5	8.810	82	3887	0.061	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	9127	0.054	ng	0.00
14) 2,4,6-Tribromophenol	15.804	330	113	0.007	ng	0.01
15) 2-Fluorobiphenyl	12.924	172	13423	0.060	ng	0.00
27) Fluoranthene-d10	19.093	212	20144	0.062	ng	0.00
31) Terphenyl-d14	19.703	244	20373	0.088	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.290	88	11544	1.128	ng	# 86

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

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