

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100621\
 Data File : BN016788.D
 Acq On : 06 Oct 2021 20:29
 Operator : CG/JU
 Sample : M3915-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D2-20210921

Manual Integrations
 APPROVED

mohammad
 10/7/2021 8:51:34 AM

Quant Time: Oct 07 00:59:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 06 14:22:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.633	152	19479	0.400	ng	0.00
7) Naphthalene-d8	10.407	136	90283	0.400	ng	# 0.00
13) Acenaphthene-d10	14.269	164	51113	0.400	ng	0.00
19) Phenanthrene-d10	17.018	188	101532	0.400	ng	0.00
29) Chrysene-d12	21.227	240	104010	0.400	ng	0.00
35) Perylene-d12	23.481	264	110669	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.264	112	6195	0.125	ng	0.02
5) Phenol-d6	6.822	99	3812	0.069	ng	0.00
8) Nitrobenzene-d5	8.772	82	15935	0.252	ng	0.00
11) 2-Methylnaphthalene-d10	11.998	152	39870	0.296	ng	0.00
14) 2,4,6-Tribromophenol	15.766	330	7973	0.388	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	51840	0.254	ng	0.00
27) Fluoranthene-d10	19.058	212	115876	0.417	ng	0.00
31) Terphenyl-d14	19.674	244	92658	0.398	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.253	88	24001m	2.640	ng	Qvalue
6) bis(2-Chloroethyl)ether	7.071	93	1287	0.024	ng	98
34) Bis(2-ethylhexyl)phtha...	21.164	149	42077	0.215	ng	100

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

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