

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
 Data File : BN016692.D
 Acq On : 03 Oct 2021 05:12
 Operator : CG/JU
 Sample : M3915-02
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D-20210921

Manual Integrations
 APPROVED

mohammad
 10/4/2021 11:49:03 AM

Quant Time: Oct 04 03:36:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 13:35:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	31848	0.40	ng	0.00
7) Naphthalene-d8	10.407	136	144338	0.40	ng	# 0.00
13) Acenaphthene-d10	14.269	164	76558	0.40	ng	0.00
19) Phenanthrene-d10	17.017	188	143614	0.40	ng	#-0.01
29) Chrysene-d12	21.227	240	130695	0.40	ng	#-0.01
35) Perylene-d12	23.481	264	128159	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.273	112	9624	0.12	ng	0.02
5) Phenol-d6	6.822	99	5619	0.06	ng	0.00
8) Nitrobenzene-d5	8.784	82	22675	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	65638	0.31	ng	0.00
14) 2,4,6-Tribromophenol	15.766	330	9773	0.28	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	84296	0.27	ng	-0.01
27) Fluoranthene-d10	19.063	212	162139	0.42	ng	0.00
31) Terphenyl-d14	19.678	244	124898	0.44	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.253	88	89145m	6.08	ng	Qvalue
6) bis(2-Chloroethyl)ether	7.080	93	4010	0.05	ng	97
34) Bis(2-ethylhexyl)phtha...	21.164	149	25928	0.16	ng	99

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

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