

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
 Data File : BN016660.D
 Acq On : 02 Oct 2021 07:33
 Operator : CG/JU
 Sample : M3829-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE125D2-20210917

Manual Integrations
 APPROVED

mohammad
 10/4/2021 11:48:21 AM

Quant Time: Oct 04 03:32:49 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	7.633	152	32644	0.40	ng	0.00
7) Naphthalene-d8	10.407	136	148411	0.40	ng	# 0.00
13) Acenaphthene-d10	14.269	164	78044	0.40	ng	0.00
19) Phenanthrene-d10	17.018	188	150856	0.40	ng	#-0.01
29) Chrysene-d12	21.227	240	132704	0.40	ng	#-0.01
35) Perylene-d12	23.481	264	135451	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.263	112	8062	0.10	ng	0.00
5) Phenol-d6	6.822	99	4693	0.05	ng	0.00
8) Nitrobenzene-d5	8.784	82	23101	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	62990	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.766	330	10200	0.29	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	85510	0.27	ng	-0.01
27) Fluoranthene-d10	19.063	212	151865	0.38	ng	0.00
31) Terphenyl-d14	19.679	244	124505	0.43	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.263	88	116515	7.75	ng	# 86
25) Phenanthrene	17.066	178	29200m	0.06	ng	
34) Bis(2-ethylhexyl)phtha...	21.164	149	20216	0.12	ng	99

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

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