

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092521\
 Data File : BN016514.D
 Acq On : 26 Sep 2021 17:01
 Operator : CG/JU
 Sample : M3833-15
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE122D1-20210920

Manual Integrations
 APPROVED

mohammad
 9/28/2021 1:37:46 PM

Quant Time: Sep 27 05:31:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 03:35:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.661	152	12294	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	42033	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	9776	0.40	ng	# 0.00
19) Phenanthrene-d10	17.042	188	30179	0.40	ng	0.00
29) Chrysene-d12	21.249	240	31289	0.40	ng	# 0.00
35) Perylene-d12	23.515	264	27753	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.291	112	1395	0.05	ng	0.00
5) Phenol-d6	6.849	99	1600	0.05	ng	0.00
8) Nitrobenzene-d5	8.810	82	11780	0.40	ng	0.00
11) 2-Methylnaphthalene-d10	12.025	152	10500	0.17	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	1132	0.44	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	17537	0.43	ng	-0.01
27) Fluoranthene-d10	19.088	212	40647	0.49	ng	0.00
31) Terphenyl-d14	19.698	244	31967	0.44	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.281	88	3938m	0.75	ng	Qvalue
25) Phenanthrene	17.090	178	2885	0.03	ng	# 89
34) Bis(2-ethylhexyl)phtha...	21.185	149	3329	0.13	ng	96

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

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