

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030375.D
 Acq On : 10 Jun 2021 15:33
 Operator : CG/JU
 Sample : M2613-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(94-98)

Manual Integrations
 APPROVED

mohammad
 6/11/2021 2:11:52 PM

Quant Time: Jun 10 16:05:35 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	185850	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	739714	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	523943	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1269060	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1358664	20.000	ng	0.00
86) Perylene-d12	23.644	264	1222912	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1515025	132.522	ng	0.00
7) Phenol-d6	7.010	99	1794815	108.449	ng	0.00
23) Nitrobenzene-d5	8.998	82	1237894	81.528	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	1004123	152.431	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	3121424	78.346	ng	0.00
79) Terphenyl-d14	19.815	244	6503520	84.076	ng	0.00
Target Compounds						
32) Benzoic acid	9.898	122	13322m	6.353	ng	Qvalue
84) Bis(2-ethylhexyl)phtha...	21.274	149	6193	4.362	ng	# 87

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

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