

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060822\
 Data File : BM035441.D
 Acq On : 08 Jun 2022 22:06
 Operator : CG/JU
 Sample : N3202-03DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-27DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/09/2022
 Supervised By : Jagrut Upadhyay 06/09/2022

Quant Time: Jun 09 00:51:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 14:28:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.839	152	220991	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	823142	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	526199	20.000	ng	-0.01
64) Phenanthrene-d10	17.221	188	1089887	20.000	ng	0.00
76) Chrysene-d12	21.409	240	1019186	20.000	ng	-0.01
86) Perylene-d12	23.762	264	1101056	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	107461	10.218	ng	0.00
7) Phenol-d6	7.034	99	77997	5.429	ng	0.00
23) Nitrobenzene-d5	8.998	82	179100	13.816	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	143614	17.010	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	561453	15.325	ng	-0.01
79) Terphenyl-d14	19.850	244	782405	15.294	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	8.633	107	41657	2.843	ng	93
27) 2,4-Dimethylphenol	9.857	122	142505m	13.985	ng	
31) Naphthalene	10.686	128	1140845	28.604	ng	100
37) 2-Methylnaphthalene	12.298	142	199909	6.643	ng	96
38) 1-Methylnaphthalene	12.521	142	97340	3.300	ng	# 98

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

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