

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060722\
 Data File : BM035420.D
 Acq On : 07 Jun 2022 20:20
 Operator : CG/JU
 Sample : N3202-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-14

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 06/08/2022
 Supervised By :Jagrut Upadhyay 06/09/2022

Quant Time: Jun 08 04:08:52 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	99648	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	368896	20.000	ng	# 0.00
39) Acenaphthene-d10	14.480	164	241063	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	509046	20.000	ng	# 0.00
76) Chrysene-d12	21.415	240	549811	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	607492	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	282818	59.639	ng	0.00
7) Phenol-d6	7.028	99	203291	31.384	ng	0.00
23) Nitrobenzene-d5	9.004	82	477531	82.195	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	511259	132.184	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	1475272	87.898	ng	0.00
79) Terphenyl-d14	19.856	244	2371058	85.917	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	9.839	122	11420m	2.501	ng	
31) Naphthalene	10.692	128	1322465	73.988	ng	99
37) 2-Methylnaphthalene	12.304	142	89856	6.662	ng	95
38) 1-Methylnaphthalene	12.521	142	187255	14.167	ng	99

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

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