

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060822\
 Data File : BM035432.D
 Acq On : 08 Jun 2022 16:32
 Operator : CG/JU
 Sample : N3189-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20220428-SPLP-AMENDED-COMPOSITE-H

Manual Integrations
 APPROVED

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

Quant Time: Jun 09 00:50:44 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	194888	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	732442	20.000	ng	# 0.00
39) Acenaphthene-d10	14.480	164	480389	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	953525	20.000	ng	0.00
76) Chrysene-d12	21.415	240	895859	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	989939	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	646174	69.672	ng	0.00
7) Phenol-d6	7.022	99	489876	38.668	ng	0.00
23) Nitrobenzene-d5	9.004	82	999250	86.626	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	933424	121.103	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	3005423	89.856	ng	0.00
79) Terphenyl-d14	19.856	244	4416396	98.215	ng	0.00

Target Compounds Qvalue

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

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