

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040422\
 Data File : BF127638.D
 Acq On : 04 Apr 2022 18:08
 Operator : CG\JU
 Sample : N2249-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2D

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/06/2022
 Supervised By : Jagrut Upadhyay 04/15/2022

Quant Time: Apr 04 21:51:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 29 07:13:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.840	152	92220	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	328107	20.000	ng	# 0.00
39) Acenaphthene-d10	9.869	164	197067	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	365987	20.000	ng	0.00
76) Chrysene-d12	14.016	240	266634	20.000	ng	0.00
86) Perylene-d12	15.498	264	203533	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	234382	40.569	ng	0.02
7) Phenol-d6	6.493	99	327521	41.473	ng	0.01
23) Nitrobenzene-d5	7.410	82	651331	84.855	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	281521	134.104	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1104523	81.843	ng	0.00
79) Terphenyl-d14	12.951	244	1232694	75.682	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	7.263	107	45053	7.331	ng	# 60
27) 2,4-Dimethylphenol	7.793	122	141119m	30.557	ng	
31) Naphthalene	8.145	128	2014812	115.663	ng	97
32) Benzoic acid	7.922	122	16364	11.432	ng	# 78
37) 2-Methylnaphthalene	8.828	142	414997	37.050	ng	100
38) 1-Methylnaphthalene	8.928	142	380908	34.391	ng	99

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

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