

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040422\
 Data File : BF127637.D
 Acq On : 04 Apr 2022 17:39
 Operator : CG\JU
 Sample : N2249-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-101R

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/05/2022
 Supervised By : Jagrut Upadhyay 04/05/2022

Quant Time: Apr 04 21:51:12 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	6.839	152	92480	20.000	ng	0.00
21) Naphthalene-d8	8.122	136	323147	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	200324	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	368280	20.000	ng	0.00
76) Chrysene-d12	14.015	240	280766	20.000	ng	0.00
86) Perylene-d12	15.498	264	210382	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	80714m	13.931	ng	0.04
7) Phenol-d6	6.492	99	299175	37.777	ng	0.01
23) Nitrobenzene-d5	7.410	82	642401	84.977	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	275291	129.005	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1048322	76.416	ng	0.00
79) Terphenyl-d14	12.951	244	1294452	75.474	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	7.263	107	50793	8.242	ng	# 59
27) 2,4-Dimethylphenol	7.798	122	235086	51.686	ng	# 94
31) Naphthalene	8.151	128	3487465	203.275	ng	97
32) Benzoic acid	7.922	122	67541	28.628	ng	# 34
37) 2-Methylnaphthalene	8.833	142	714620	64.779	ng	100
38) 1-Methylnaphthalene	8.928	142	410770	37.656	ng	98

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

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