

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060721\
 Data File : BF124155.D
 Acq On : 7 Jun 2021 14:01
 Operator : JU/CG
 Sample : M2583-22
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 54-57

Manual Integrations
 APPROVED

mohammad
 6/8/2021 3:36:27 PM

Quant Time: Jun 07 14:22:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 07 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	37298	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	151816	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	85115	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	128536	20.000	ng	0.00
76) Chrysene-d12	14.039	240	85759	20.000	ng	0.00
86) Perylene-d12	15.527	264	86484	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	202819	81.857	ng	0.01
7) Phenol-d6	6.522	99	171043	51.358	ng	0.02
23) Nitrobenzene-d5	7.434	82	324353	84.736	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	135503	112.369	ng	0.01
45) 2-Fluorobiphenyl	9.234	172	592313	87.569	ng	0.01
79) Terphenyl-d14	12.980	244	362774	58.078	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.534	94	171401	47.622	ng	88
20) 3+4-Methylphenols	7.292	107	616011m	207.438	ng	
22) Acetophenone	7.275	105	12228m	2.825	ng	
27) 2,4-Dimethylphenol	7.892	122	24053	9.944	ng	93
36) 4-Chloro-3-methylphenol	8.745	107	193022	63.034	ng	# 85
50) Dimethylphthalate	9.639	163	58451	8.300	ng	# 98
60) Diethylphthalate	10.339	149	21413	3.013	ng	# 96
70) Pentachlorophenol	11.210	266	7564	11.429	ng	91
84) Bis(2-ethylhexyl)phtha...	14.016	149	23389	5.886	ng	94

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

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