

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020824\
 Data File : BF137340.D
 Acq On : 08 Feb 2024 16:24
 Operator : CG\JU
 Sample : P1372-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CG-MW-17-ch5

Quant Time: Feb 08 16:50:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 04:56:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.786	152	102010	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	386372	20.000	ng	0.00
39) Acenaphthene-d10	9.816	164	213972	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	364377	20.000	ng	0.00
76) Chrysene-d12	13.962	240	221045	20.000	ng	-0.01
86) Perylene-d12	15.439	264	248068	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	391631	63.349	ng	0.00
7) Phenol-d6	6.416	99	307871	34.896	ng	-0.01
23) Nitrobenzene-d5	7.351	82	742809	86.074	ng	0.00
42) 2,4,6-Tribromophenol	10.610	330	285652	118.209	ng	-0.01
45) 2-Fluorobiphenyl	9.145	172	1255209	80.641	ng	0.00
79) Terphenyl-d14	12.910	244	1155625	71.072	ng	-0.01
Target Compounds						
						Qvalue
10) Phenol	6.428	94	85004	8.632	ng	94
20) 3+4-Methylphenols	7.204	107	1396660	191.048	ng	# 57
31) Naphthalene	8.098	128	3514418	168.252	ng	97
38) 1-Methylnaphthalene	8.875	142	609414	46.854	ng	98
49) Acenaphthylene	9.675	152	335108	16.614	ng	99
52) Acenaphthene	9.851	154	207510	17.505	ng	99
73) Carbazole	11.545	167	36087	2.172	ng	99

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

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