

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050123\
 Data File : BF133167.D
 Acq On : 01 May 2023 21:13
 Operator : CG\JU
 Sample : 02558-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS012-0006-01

Quant Time: May 02 05:23:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 01 15:02:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	212779	20.000	ng	0.00
21) Naphthalene-d8	8.016	136	698927	20.000	ng	0.00
39) Acenaphthene-d10	9.769	164	365704	20.000	ng	0.00
64) Phenanthrene-d10	11.251	188	605822	20.000	ng	0.00
76) Chrysene-d12	13.886	240	307015	20.000	ng	# 0.00
86) Perylene-d12	15.315	264	419020	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	1090093	77.390	ng	0.00
7) Phenol-d6	6.369	99	1331358	76.151	ng	0.00
23) Nitrobenzene-d5	7.292	82	655502	50.623	ng	-0.01
42) 2,4,6-Tribromophenol	10.557	330	266703	72.516	ng	0.00
45) 2-Fluorobiphenyl	9.092	172	1047383	47.915	ng	0.00
79) Terphenyl-d14	12.839	244	778543	43.735	ng	0.00
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	7.139	107	134879	8.746	ng	# 55
31) Naphthalene	8.034	128	358994	10.273	ng	# 97
37) 2-Methylnaphthalene	8.728	142	699944	29.298	ng	99
38) 1-Methylnaphthalene	8.828	142	484340	21.672	ng	100
46) 1,1'-Biphenyl	9.192	154	82136	2.832	ng	97
58) Fluorene	10.316	166	86229	3.799	ng	# 88
71) Phenanthrene	11.274	178	229665	7.053	ng	96

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

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