

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061322\
 Data File : BF128785.D
 Acq On : 13 Jun 2022 15:41
 Operator : CG\JU
 Sample : N3286-01DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1RDL

Quant Time: Jun 14 01:55:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 08 14:28:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.804	152	90030	20.000	ng	0.00
21) Naphthalene-d8	8.086	136	348777	20.000	ng	-0.01
39) Acenaphthene-d10	9.845	164	201553	20.000	ng	-0.01
64) Phenanthrene-d10	11.333	188	386273	20.000	ng	0.00
76) Chrysene-d12	13.980	240	297708	20.000	ng	# 0.00
86) Perylene-d12	15.445	264	210956	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	30719	5.435	ng	0.00
7) Phenol-d6	6.463	99	28045	4.065	ng	-0.02
23) Nitrobenzene-d5	7.363	82	56772	7.579	ng	-0.02
42) 2,4,6-Tribromophenol	10.633	330	26557	11.152	ng	-0.02
45) 2-Fluorobiphenyl	9.163	172	118963	8.268	ng	-0.01
79) Terphenyl-d14	12.915	244	144303	8.423	ng	-0.02
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	7.227	107	19435	3.185	ng	# 75
27) 2,4-Dimethylphenol	7.751	122	10223	2.215	ng	# 82
31) Naphthalene	8.110	128	755184	42.190	ng	99
37) 2-Methylnaphthalene	8.798	142	85957	7.550	ng	99
38) 1-Methylnaphthalene	8.898	142	49529	4.503	ng	96

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

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