

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061322\
 Data File : BF128787.D
 Acq On : 13 Jun 2022 16:42
 Operator : CG\JU
 Sample : N3286-03DL 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2DDL

Quant Time: Jun 14 01:56:06 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	96567	20.000	ng	0.00
21) Naphthalene-d8	8.087	136	364438	20.000	ng	#-0.01
39) Acenaphthene-d10	9.845	164	207346	20.000	ng	-0.01
64) Phenanthrene-d10	11.333	188	384211	20.000	ng	0.00
76) Chrysene-d12	13.974	240	268529	20.000	ng	#-0.01
86) Perylene-d12	15.439	264	184562	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	141984	23.420	ng	0.00
7) Phenol-d6	6.463	99	115706	15.635	ng	-0.02
23) Nitrobenzene-d5	7.369	82	299366	38.245	ng	-0.01
42) 2,4,6-Tribromophenol	10.639	330	145246	59.287	ng	-0.01
45) 2-Fluorobiphenyl	9.163	172	610381	41.236	ng	-0.01
79) Terphenyl-d14	12.922	244	674019	43.615	ng	-0.01
Target Compounds						
27) 2,4-Dimethylphenol	7.769	122	49486	10.260	ng	# 92
31) Naphthalene	8.110	128	682767	36.505	ng	99
37) 2-Methylnaphthalene	8.798	142	102125	8.585	ng	97
38) 1-Methylnaphthalene	8.898	142	115413	10.041	ng	98

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

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