

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052722\
 Data File : BF128507.D
 Acq On : 27 May 2022 13:39
 Operator : CG\JU
 Sample : N3056-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-02

Quant Time: May 27 23:18:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	194710	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	694264	20.000	ng	0.00
39) Acenaphthene-d10	9.880	164	425947	20.000	ng	0.00
64) Phenanthrene-d10	11.368	188	783057	20.000	ng	0.00
76) Chrysene-d12	14.004	240	474507	20.000	ng	0.00
86) Perylene-d12	15.468	264	393683	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	576123	49.011	ng	0.00
7) Phenol-d6	6.469	99	644027	42.970	ng	0.00
23) Nitrobenzene-d5	7.416	82	1166885	96.959	ng	0.00
42) 2,4,6-Tribromophenol	10.674	330	600666	148.155	ng	0.00
45) 2-Fluorobiphenyl	9.204	172	2232276	85.143	ng	0.00
79) Terphenyl-d14	12.957	244	2414583	98.310	ng	0.00
Target Compounds						
						Qvalue
27) 2,4-Dimethylphenol	7.781	122	156191	15.802	ng	97
31) Naphthalene	8.157	128	5761041	169.543	ng	94
37) 2-Methylnaphthalene	8.839	142	1144298	53.352	ng	99
38) 1-Methylnaphthalene	8.939	142	956363	45.846	ng	98
58) Fluorene	10.427	166	60423	2.542	ng	98
71) Phenanthrene	11.386	178	109403	2.949	ng	98
73) Carbazole	11.592	167	222603	6.269	ng	99

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

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