

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061722\
 Data File : BF128857.D
 Acq On : 17 Jun 2022 16:39
 Operator : CG\JU
 Sample : N3373-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1E

Quant Time: Jun 18 05:06:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 14 14:53:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	82988	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	319335	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	147477	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	217458	20.000	ng	# 0.00
76) Chrysene-d12	13.957	240	195070	20.000	ng	# 0.00
86) Perylene-d12	15.409	264	190226	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	653893	125.505	ng	0.00
7) Phenol-d6	6.439	99	849165	133.520	ng	0.00
23) Nitrobenzene-d5	7.351	82	601070	87.635	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	164294	94.286	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	988172	93.861	ng	0.00
79) Terphenyl-d14	12.904	244	796738	70.971	ng	0.00
Target Compounds						
						Qvalue
55) Dibenzofuran	10.022	168	30850	2.550	ng	# 91
71) Phenanthrene	11.357	178	3691938	338.485	ng	99
72) Anthracene	11.398	178	455753	42.216	ng	98
73) Carbazole	11.557	167	742949	74.094	ng	98
75) Fluoranthene	12.580	202	8866788	777.698	ng	95
78) Pyrene	12.804	202	5337596	367.777	ng	99
81) Benzo(a)anthracene	13.945	228	172244	14.166	ng	98
83) Chrysene	13.986	228	205447	17.722	ng	98

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

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