

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112822\
 Data File : BF131440.D
 Acq On : 28 Nov 2022 21:00
 Operator : CG\JU
 Sample : N5742-10DL 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP111822DL

Quant Time: Nov 29 00:59:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 00:18:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	141794	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	497325	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	253725	20.000	ng	0.00
64) Phenanthrene-d10	11.504	188	396651	20.000	ng	0.00
76) Chrysene-d12	14.157	240	339321	20.000	ng	#-0.01
86) Perylene-d12	15.698	264	331924	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.557	112	48317	6.485	ng	0.00
7) Phenol-d6	6.563	99	54807	5.766	ng	-0.01
23) Nitrobenzene-d5	7.516	82	280656	28.447	ng	-0.01
42) 2,4,6-Tribromophenol	10.798	330	60889	28.525	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	584299	33.503	ng	0.00
79) Terphenyl-d14	13.092	244	439652	21.188	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	8.269	128	2124739	80.172	ng	99
37) 2-Methylnaphthalene	8.963	142	728738	40.583	ng	100
38) 1-Methylnaphthalene	9.057	142	487617	28.003	ng	99
46) 1,1'-Biphenyl	9.422	154	49155	2.539	ng	99

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

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