

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100422\
 Data File : BF130522.D
 Acq On : 04 Oct 2022 17:14
 Operator : CG\JU
 Sample : N4941-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6R

Quant Time: Oct 05 02:19:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.634	152	76361	20.000	ng	-0.01
21) Naphthalene-d8	7.910	136	286300	20.000	ng	-0.02
39) Acenaphthene-d10	9.663	164	132924	20.000	ng	-0.02
64) Phenanthrene-d10	11.145	188	193048	20.000	ng	-0.01
76) Chrysene-d12	13.774	240	143741	20.000	ng	-0.02
86) Perylene-d12	15.162	264	175575	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	192161	43.647	ng	-0.01
7) Phenol-d6	6.269	99	157090	28.517	ng	-0.02
23) Nitrobenzene-d5	7.198	82	494094	88.168	ng	-0.02
42) 2,4,6-Tribromophenol	10.451	330	110998	87.148	ng	-0.02
45) 2-Fluorobiphenyl	8.992	172	850573	93.181	ng	-0.02
79) Terphenyl-d14	12.727	244	612169	70.547	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	7.933	128	90382	6.128	ng	100
52) Acenaphthene	9.692	154	125749	15.895	ng	99
55) Dibenzofuran	9.869	168	64969	5.904	ng	96

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

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