

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042823\
 Data File : BF133143.D
 Acq On : 29 Apr 2023 02:40
 Operator : CG\JU
 Sample : 02522-03DL 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12DL

Quant Time: May 01 03:11:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 01 03:03:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.734	152	212031	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	874371	20.000	ng	0.00
39) Acenaphthene-d10	9.769	164	442161	20.000	ng	0.00
64) Phenanthrene-d10	11.251	188	801778	20.000	ng	0.00
76) Chrysene-d12	13.880	240	580600	20.000	ng	-0.01
86) Perylene-d12	15.304	264	460102	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	90572	6.453	ng	0.00
7) Phenol-d6	6.363	99	65065	3.735	ng	-0.02
23) Nitrobenzene-d5	7.292	82	145407	8.976	ng	-0.01
42) 2,4,6-Tribromophenol	10.551	330	62118	13.969	ng	-0.01
45) 2-Fluorobiphenyl	9.092	172	295323	11.174	ng	0.00
79) Terphenyl-d14	12.833	244	345764	10.271	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.051	128	4052701	92.706	ng	98
37) 2-Methylnaphthalene	8.733	142	720045	24.092	ng	100
38) 1-Methylnaphthalene	8.833	142	1242096	44.425	ng	99
46) 1,1'-Biphenyl	9.192	154	238490	6.802	ng	97
52) Acenaphthene	9.798	154	230267	8.385	ng	99
58) Fluorene	10.310	166	100330	3.655	ng	100
71) Phenanthrene	11.269	178	207581	4.817	ng	98

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

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