

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101123\
 Data File : BF135716.D
 Acq On : 11 Oct 2023 16:42
 Operator : CG\JU
 Sample : 04656-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-103S

Quant Time: Oct 12 00:48:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 02:43:17 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.739	152	72262	20.000	ng	0.00
21) Naphthalene-d8	8.016	136	271092	20.000	ng	0.00
39) Acenaphthene-d10	9.775	164	135558	20.000	ng	0.00
64) Phenanthrene-d10	11.269	188	237020	20.000	ng	0.00
76) Chrysene-d12	13.933	240	146441	20.000	ng	0.00
86) Perylene-d12	15.404	264	186931	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	249349	56.123	ng	0.00
7) Phenol-d6	6.392	99	191969	33.980	ng	0.00
23) Nitrobenzene-d5	7.310	82	478301	95.856	ng	0.00
42) 2,4,6-Tribromophenol	10.575	330	196636	145.968	ng	0.00
45) 2-Fluorobiphenyl	9.098	172	872816	97.142	ng	0.00
79) Terphenyl-d14	12.868	244	576705	61.049	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.039	128	50885	3.863	ng	99
52) Acenaphthene	9.804	154	35719	4.813	ng	99
71) Phenanthrene	11.292	178	23417	2.089	ng	98
75) Fluoranthene	12.498	202	28186	2.413	ng	96

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

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