

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011524\
 Data File : BF137068.D
 Acq On : 15 Jan 2024 17:38
 Operator : CG\JU
 Sample : P1135-01DL 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 360DL

Quant Time: Jan 15 18:20:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	42075	20.000	ng	-0.01
21) Naphthalene-d8	8.063	136	154114	20.000	ng	# 0.00
39) Acenaphthene-d10	9.816	164	73288	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	144842	20.000	ng	0.00
76) Chrysene-d12	13.969	240	143959	20.000	ng	0.00
86) Perylene-d12	15.463	264	136899	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	116595	43.237	ng	0.00
7) Phenol-d6	6.434	99	137845	39.450	ng	0.00
23) Nitrobenzene-d5	7.345	82	81770	27.150	ng	-0.01
42) 2,4,6-Tribromophenol	10.610	330	36289	42.044	ng	-0.01
45) 2-Fluorobiphenyl	9.134	172	155109	28.466	ng	-0.02
79) Terphenyl-d14	12.904	244	184682	17.928	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	8.081	128	427939	50.544	ng	100
37) 2-Methylnaphthalene	8.775	142	85306	15.656	ng	100
38) 1-Methylnaphthalene	8.869	142	49359	9.233	ng	96
46) 1,1'-Biphenyl	9.239	154	23211	3.774	ng	96
52) Acenaphthene	9.851	154	66987	15.122	ng	98
55) Dibenzofuran	10.022	168	95260	14.186	ng	95
58) Fluorene	10.369	166	67879	12.740	ng	98
71) Phenanthrene	11.333	178	263237	35.420	ng	99
72) Anthracene	11.386	178	23386	3.107	ng	99
73) Carbazole	11.551	167	30846	4.347	ng	99
75) Fluoranthene	12.533	202	130169	16.374	ng	99
78) Pyrene	12.763	202	89805	6.355	ng	97

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

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