

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010224\
 Data File : BF136874.D
 Acq On : 02 Jan 2024 16:38
 Operator : CG\JU
 Sample : 06038-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T-401

Quant Time: Jan 03 01:36:47 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	56443	20.000	ng	-0.01
21) Naphthalene-d8	8.063	136	203898	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	100367	20.000	ng	-0.02
64) Phenanthrene-d10	11.304	188	162976	20.000	ng	-0.01
76) Chrysene-d12	13.957	240	105748	20.000	ng	-0.01
86) Perylene-d12	15.439	264	132637	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	374061	103.404	ng	0.00
7) Phenol-d6	6.428	99	429958	91.727	ng	-0.01
23) Nitrobenzene-d5	7.345	82	301714	75.719	ng	-0.01
42) 2,4,6-Tribromophenol	10.610	330	130048	110.020	ng	-0.01
45) 2-Fluorobiphenyl	9.133	172	586857	78.643	ng	-0.02
79) Terphenyl-d14	12.898	244	531229	70.202	ng	-0.02
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	7.198	107	24484	5.976	ng	# 51
31) Naphthalene	8.098	128	3863433	344.900	ng	99
37) 2-Methylnaphthalene	8.775	142	441170	61.197	ng	100
38) 1-Methylnaphthalene	8.869	142	317847	44.940	ng	99
46) 1,1'-Biphenyl	9.233	154	54651	6.489	ng	97
52) Acenaphthene	9.845	154	80803	13.320	ng	96
58) Fluorene	10.363	166	43252	5.928	ng	97
71) Phenanthrene	11.327	178	83741	10.014	ng	98

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

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