

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022323\
 Data File : BF132525.D
 Acq On : 23 Feb 2023 17:41
 Operator : CG\JU
 Sample : 01635-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-294

Quant Time: Feb 24 00:38:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 23 00:59:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.007	152	188511	20.000	ng	0.00
21) Naphthalene-d8	8.295	136	814328	20.000	ng	0.00
39) Acenaphthene-d10	10.054	164	464222	20.000	ng	0.00
64) Phenanthrene-d10	11.554	188	865693	20.000	ng	0.00
76) Chrysene-d12	14.207	240	439737	20.000	ng	-0.01
86) Perylene-d12	15.760	264	444507	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	1046019	90.056	ng	0.00
7) Phenol-d6	6.637	99	1403015	92.554	ng	0.00
23) Nitrobenzene-d5	7.572	82	941337	54.844	ng	0.00
42) 2,4,6-Tribromophenol	10.854	330	357565	71.322	ng	0.00
45) 2-Fluorobiphenyl	9.372	172	1535361	50.360	ng	0.00
79) Terphenyl-d14	13.142	244	1652292	63.529	ng	0.00
Target Compounds						
						Qvalue
9) Benzaldehyde	6.560	77	332272	40.622	ng	98
10) Phenol	6.649	94	112224	6.457	ng	97
22) Acetophenone	7.407	105	55900	2.722	ng	# 86
32) Benzoic acid	8.137	122	988988	94.170	ng	92
46) 1,1'-Biphenyl	9.472	154	560767	15.960	ng	98

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

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