

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022823\
 Data File : BF132578.D
 Acq On : 28 Feb 2023 18:25
 Operator : CG\JU
 Sample : 01701-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-13-GW01

Quant Time: Mar 01 02:29:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 24 23:17:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.004	152	239085	20.000	ng	0.00
21) Naphthalene-d8	8.287	136	866856	20.000	ng	0.00
39) Acenaphthene-d10	10.051	164	466608	20.000	ng	0.00
64) Phenanthrene-d10	11.545	188	925506	20.000	ng	0.00
76) Chrysene-d12	14.198	240	608526	20.000	ng	# 0.00
86) Perylene-d12	15.751	264	458594	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.634	112	1081768	73.433	ng	0.00
7) Phenol-d6	6.622	99	1046735	54.445	ng	-0.01
23) Nitrobenzene-d5	7.569	82	1691024	92.552	ng	0.00
42) 2,4,6-Tribromophenol	10.851	330	758411	150.503	ng	0.00
45) 2-Fluorobiphenyl	9.369	172	2630594	85.843	ng	0.00
79) Terphenyl-d14	13.139	244	3463450	96.229	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	8.122	122	12376	4.898	ng	91
54) 2,4-Dinitrophenol	10.104	184	4491	2.745	ng	91
56) 4-Nitrophenol	10.163	139	13023	2.018	ng	92
70) Pentachlorophenol	11.345	266	15170	2.315	ng	96
92) Benzo(g,h,i)perylene	17.833	276	31566	2.717	ng	94

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

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