

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022723\
 Data File : BF132563.D
 Acq On : 27 Feb 2023 20:53
 Operator : CG\JU
 Sample : 01701-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-19-SB01

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/28/2023
 Supervised By :Jagrut Upadhyay 02/28/2023

Quant Time: Feb 28 03:18:34 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	7.001	152	147002	20.000	ng	0.00
21) Naphthalene-d8	8.289	136	520814	20.000	ng	0.00
39) Acenaphthene-d10	10.048	164	255862	20.000	ng	0.00
64) Phenanthrene-d10	11.548	188	412361	20.000	ng	# 0.00
76) Chrysene-d12	14.207	240	315404	20.000	ng	0.00
86) Perylene-d12	15.760	264	238372	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.631	112	693904	76.610	ng	0.00
7) Phenol-d6	6.625	99	896787	75.864	ng	-0.01
23) Nitrobenzene-d5	7.566	82	599310	54.595	ng	0.00
42) 2,4,6-Tribromophenol	10.842	330	189746	68.669	ng	0.00
45) 2-Fluorobiphenyl	9.366	172	988978	58.855	ng	0.00
79) Terphenyl-d14	13.136	244	866467	46.448	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.307	128	127893	4.797	ng	98
37) 2-Methylnaphthalene	9.001	142	120183	6.614	ng	96
38) 1-Methylnaphthalene	9.101	142	103206	6.078	ng	100
55) Dibenzofuran	10.254	168	60471	2.673	ng	# 98
71) Phenanthrene	11.572	178	420146	19.668	ng	98
72) Anthracene	11.619	178	90782	4.127	ng	99
75) Fluoranthene	12.766	202	521099	22.314	ng	99
78) Pyrene	13.001	202	510413	17.799	ng	99
81) Benzo(a)anthracene	14.195	228	276719	11.726	ng	94
83) Chrysene	14.230	228	298327	13.518	ng	96
87) Indeno(1,2,3-cd)pyrene	17.354	276	92246	5.906	ng	95
88) Benzo(b)fluoranthene	15.295	252	301557m	18.877	ng	
89) Benzo(k)fluoranthene	15.318	252	95810m	6.128	ng	
90) Benzo(a)pyrene	15.689	252	172293	11.524	ng	96
92) Benzo(g,h,i)perylene	17.842	276	92013	7.797	ng	99

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

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