

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101923\
 Data File : BF135898.D
 Acq On : 19 Oct 2023 21:22
 Operator : CG\JU
 Sample : 04704-06DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-4(5-10)DL

Quant Time: Oct 20 02:05:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 17 01:22:00 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/20/2023
 Supervised By :mohammad ahmed 10/20/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.734	152	67929	20.000	ng	-0.01
21) Naphthalene-d8	8.016	136	221544	20.000	ng	0.00
39) Acenaphthene-d10	9.769	164	107193	20.000	ng	-0.02
64) Phenanthrene-d10	11.269	188	155264	20.000	ng	-0.01
76) Chrysene-d12	13.933	240	104560	20.000	ng	-0.02
86) Perylene-d12	15.415	264	122612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	64318	15.223	ng	0.00
7) Phenol-d6	6.398	99	108347	20.552	ng	0.00
23) Nitrobenzene-d5	7.304	82	50420	12.517	ng	-0.01
42) 2,4,6-Tribromophenol	10.574	330	12108	11.771	ng	-0.01
45) 2-Fluorobiphenyl	9.092	172	126510	18.453	ng	-0.01
79) Terphenyl-d14	12.863	244	109152	19.332	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.033	128	121222	11.218	ng	99
37) 2-Methylnaphthalene	8.733	142	31624	4.442	ng	90
38) 1-Methylnaphthalene	8.828	142	45200	6.883	ng	99
49) Acenaphthylene	9.633	152	49663	5.071	ng	99
52) Acenaphthene	9.804	154	17214	2.933	ng	93
55) Dibenzofuran	9.980	168	24567	2.774	ng	# 96
58) Fluorene	10.322	166	45620	6.579	ng	100
71) Phenanthrene	11.292	178	245518	33.009	ng	99
72) Anthracene	11.345	178	70130	9.099	ng	96
75) Fluoranthene	12.492	202	253165	30.668	ng	97
78) Pyrene	12.721	202	285164	34.320	ng	99
81) Benzo(a)anthracene	13.927	228	106090	15.447	ng	98
83) Chrysene	13.962	228	102885	15.467	ng	98
87) Indeno(1,2,3-cd)pyrene	16.921	276	30924	4.629	ng	96
88) Benzo(b)fluoranthene	14.986	252	97896m	14.266	ng	
89) Benzo(k)fluoranthene	15.009	252	28940m	4.344	ng	
90) Benzo(a)pyrene	15.357	252	89070	13.803	ng	97
92) Benzo(g,h,i)perylene	17.380	276	31480	5.596	ng	97

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

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