

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100923\
 Data File : BF135672.D
 Acq On : 10 Oct 2023 02:17
 Operator : CG\JU
 Sample : 04609-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2

Quant Time: Oct 10 04:23:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 04 17:06:20 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/11/2023
 Supervised By :mohammad ahmed 10/11/2023

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 ahmed

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	108279	20.000	ng	0.00
21) Naphthalene-d8	8.016	136	421324	20.000	ng	0.00
39) Acenaphthene-d10	9.775	164	196814	20.000	ng	0.00
64) Phenanthrene-d10	11.269	188	324409	20.000	ng	0.00
76) Chrysene-d12	13.933	240	204879	20.000	ng	0.00
86) Perylene-d12	15.404	264	190504	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	717064	107.709	ng	0.01
7) Phenol-d6	6.393	99	908137	107.278	ng	0.00
23) Nitrobenzene-d5	7.304	82	567134	73.131	ng	0.00
42) 2,4,6-Tribromophenol	10.575	330	199421	101.961	ng	0.00
45) 2-Fluorobiphenyl	9.098	172	1010194	77.439	ng	0.00
79) Terphenyl-d14	12.869	244	809688	61.264	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.040	128	75696	3.698	ng	98
52) Acenaphthene	9.804	154	77962	7.236	ng	99
55) Dibenzofuran	9.981	168	96283	5.856	ng	98
58) Fluorene	10.328	166	101114	8.000	ng	98
71) Phenanthrene	11.298	178	915729	59.683	ng	99
72) Anthracene	11.345	178	311859	19.774	ng	99
73) Carbazole	11.510	167	77255	5.381	ng	99
75) Fluoranthene	12.498	202	1132434	70.833	ng	98
78) Pyrene	12.727	202	948851	48.644	ng	100
81) Benzo(a)anthracene	13.921	228	525266	39.704	ng	98
83) Chrysene	13.963	228	428119	32.549	ng	98
87) Indeno(1,2,3-cd)pyrene	16.892	276	167441	13.405	ng	96
88) Benzo(b)fluoranthene	14.980	252	510471m	49.499	ng	
89) Benzo(k)fluoranthene	15.004	252	137470m	13.495	ng	
90) Benzo(a)pyrene	15.345	252	325093	33.056	ng	99
91) Dibenzo(a,h)anthracene	16.892	278	46885m	4.588	ng	
92) Benzo(g,h,i)perylene	17.345	276	147975	13.864	ng	96

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

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