

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090523\
 Data File : BF135109.D
 Acq On : 05 Sep 2023 23:32
 Operator : CG\JU
 Sample : 04265-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-3-083123

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/06/2023
 Supervised By :mohammad ahmed 09/07/2023

Quant Time: Sep 06 01:59:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 31 03:48:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.775	152	98297	20.000	ng	0.00
21) Naphthalene-d8	8.051	136	387731	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	169657	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	260162	20.000	ng	0.00
76) Chrysene-d12	13.957	240	175180	20.000	ng	0.00
86) Perylene-d12	15.427	264	156078	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	553596	86.991	ng	0.00
7) Phenol-d6	6.410	99	665671	83.921	ng	0.00
23) Nitrobenzene-d5	7.334	82	415926	60.266	ng	-0.01
42) 2,4,6-Tribromophenol	10.598	330	128534	71.419	ng	-0.01
45) 2-Fluorobiphenyl	9.128	172	711196	68.250	ng	-0.01
79) Terphenyl-d14	12.898	244	578438	53.885	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	8.863	142	25492	2.156	ng	99
52) Acenaphthene	9.839	154	58209	6.199	ng	99
55) Dibenzofuran	10.016	168	37750	2.787	ng	95
58) Fluorene	10.357	166	68009	6.586	ng	97
71) Phenanthrene	11.327	178	489071	36.438	ng	99
72) Anthracene	11.374	178	131698	9.604	ng	99
73) Carbazole	11.533	167	33140	2.802	ng	# 92
75) Fluoranthene	12.521	202	546915	40.303	ng	100
78) Pyrene	12.751	202	479646	28.838	ng	100
81) Benzo(a)anthracene	13.945	228	240231	20.104	ng	96
83) Chrysene	13.986	228	213313	18.164	ng	98
88) Benzo(b)fluoranthene	15.004	252	203937m	21.000	ng	
89) Benzo(k)fluoranthene	15.027	252	77388m	8.248	ng	
90) Benzo(a)pyrene	15.362	252	149852	16.610	ng	98
91) Dibenzo(a,h)anthracene	16.921	278	19979	2.346	ng	# 83
92) Benzo(g,h,i)perylene	17.356	276	73397	8.288	ng	96

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

