

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060821\
 Data File : BF124195.D
 Acq On : 8 Jun 2021 23:29
 Operator : JU/CG
 Sample : M2589-12RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 18-B10(2-4)RE

Manual Integrations
 APPROVED

mohammad
 6/9/2021 3:51:38 PM

Quant Time: Jun 09 04:10:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 07 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	39534	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	136595	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	64668	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	104000	20.000	ng	# 0.00
76) Chrysene-d12	14.045	240	89504	20.000	ng	0.01
86) Perylene-d12	15.533	264	63234	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	7248	2.760	ng	0.03
7) Phenol-d6	6.510	99	68849	19.504	ng	0.00
23) Nitrobenzene-d5	7.428	82	200460	58.205	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	9.222	172	320860	62.436	ng	0.00
79) Terphenyl-d14	12.980	244	316084	48.486	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.169	128	97928	11.991	ng	98
37) 2-Methylnaphthalene	8.857	142	30453	5.317	ng	88
38) 1-Methylnaphthalene	8.957	142	21543	4.025	ng	81
52) Acenaphthene	9.933	154	33579	7.960	ng	95
55) Dibenzofuran	10.104	168	40852	6.187	ng	95
58) Fluorene	10.451	166	31685	6.252	ng	# 97
71) Phenanthrene	11.416	178	426500	65.649	ng	97
72) Anthracene	11.469	178	93470	14.286	ng	95
73) Carbazole	11.622	167	35367	6.203	ng	95
75) Fluoranthene	12.616	202	523842	76.816	ng	97
78) Pyrene	12.845	202	475591	55.335	ng	99
81) Benzo(a)anthracene	14.033	228	204956	31.433	ng	97
83) Chrysene	14.068	228	164208	26.103	ng	95
84) Bis(2-ethylhexyl)phtha...	14.016	149	258044	62.219	ng	99
87) Indeno(1,2,3-cd)pyrene	17.033	276	53381	13.562	ng	98
88) Benzo(b)fluoranthene	15.098	252	166903m	33.978	ng	
89) Benzo(k)fluoranthene	15.121	252	46796m	10.063	ng	
90) Benzo(a)pyrene	15.468	252	110558	25.913	ng	# 93
91) Dibenzo(a,h)anthracene	17.045	278	16549	7.539	ng	# 70
92) Benzo(g,h,i)perylene	17.492	276	56613	15.566	ng	97

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

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