

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090519\
 Data File : BF116846.D
 Acq On : 5 Sep 2019 18:55
 Operator : HP/JU
 Sample : K4692-14 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-26-COMP

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:55:19 AM

Quant Time: Sep 06 14:47:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	66049	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	275529	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	143192	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	211075	20.00	ng	0.00
76) Chrysene-d12	14.02	240	148471	20.00	ng	0.02
87) Perylene-d12	15.47	264	127577	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	253298	62.13	ng	0.00
7) Phenol-d6	6.47	99	355527	66.41	ng	0.00
23) Nitrobenzene-d5	7.40	82	212772	44.08	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	69867	72.99	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	412634	48.61	ng	-0.01
79) Terphenyl-d14	12.95	244	368555	45.92	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.49	94	19127	3.227	ng	81
20) 3+4-Methylphenols	7.26	107	9436	2.083	ng	# 71
31) Naphthalene	8.15	128	165382	12.623	ng	100
37) 2-Methylnaphthalene	8.84	142	49729	5.705	ng	96
38) 1-Methylnaphthalene	8.94	142	42546m	5.170	ng	
46) 1,1'-Biphenyl	9.30	154	24727	2.276	ng	98
49) Acenaphthylene	9.75	152	490967	37.741	ng	99
50) Dimethylphthalate	9.59	163	57768	5.843	ng	98
52) Acenaphthene	9.91	154	91580	11.457	ng	97
55) Dibenzofuran	10.09	168	192529	17.086	ng	98
58) Fluorene	10.43	166	317434	36.972	ng	99
71) Phenanthrene	11.41	178	2828195	240.702	ng	97
72) Anthracene	11.45	178	413430	34.954	ng	100
73) Carbazole	11.59	167	217201	19.278	ng	99
75) Fluoranthene	12.60	202	3233692	280.043	ng	96
78) Pyrene	12.83	202	2833693	214.705	ng	97
81) Benzo(a)anthracene	14.00	228	944742m	100.540	ng	
83) Chrysene	14.04	228	1322966m	136.657	ng	
84) Bis(2-ethylhexyl)phthalate	13.98	149	26423	3.335	ng	# 92
86) Indeno(1,2,3-cd)pyrene	16.94	276	410104	52.740	ng	98
88) Benzo(b)fluoranthene	15.06	252	1260464m	163.418	ng	
89) Benzo(k)fluoranthene	15.08	252	377613m	53.694	ng	
90) Benzo(a)pyrene	15.43	252	684451	101.997	ng	99
91) Dibenzo(a,h)anthracene	16.94	278	98713m	16.806	ng	
92) Benzo(g,h,i)perylene	17.39	276	365851	61.072	ng	98

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

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