

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090519\
 Data File : BF116848.D
 Acq On : 5 Sep 2019 19:48
 Operator : HP/JU
 Sample : K4692-10 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-28-COMP

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:55:25 AM

Quant Time: Sep 06 05:30:57 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	64178	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	266904	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	136846	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	207416	20.00	ng	0.00
76) Chrysene-d12	14.00	240	154836	20.00	ng	0.00
87) Perylene-d12	15.46	264	114356	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	245031	61.85	ng	0.00
7) Phenol-d6	6.47	99	345144	66.35	ng	0.00
23) Nitrobenzene-d5	7.40	82	202100	43.23	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	62714	68.55	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	399712	49.27	ng	-0.01
79) Terphenyl-d14	12.94	244	351642	42.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	54036	5.719	ng	96
71) Phenanthrene	11.39	178	222109	19.237	ng	99
72) Anthracene	11.44	178	49767	4.282	ng	99
75) Fluoranthene	12.58	202	363899	32.070	ng	99
78) Pyrene	12.81	202	355559	25.833	ng	99
81) Benzo(a)anthracene	13.99	228	172303	17.583	ng	97
83) Chrysene	14.03	228	159812	15.829	ng	98
86) Indeno(1,2,3-cd)pyrene	16.92	276	49196	6.067	ng	97
88) Benzo(b)fluoranthene	15.04	252	153172m	22.155	ng	
89) Benzo(k)fluoranthene	15.06	252	34933m	5.542	ng	
90) Benzo(a)pyrene	15.40	252	94479	15.707	ng	100
91) Dibenzo(a,h)anthracene	16.92	278	13902	2.640	ng	# 83
92) Benzo(a,h,i)perylene	17.36	276	48964	9.119	ng	# 91

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

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