

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\
 Data File : BF115888.D
 Acq On : 1 Aug 2019 1:12
 Operator : HP/JU
 Sample : K3622-15 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-072919-B

Manual Integrations
 APPROVED

mohammad
 8/1/2019 5:07:21 PM

Quant Time: Aug 01 08:37:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	138663	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	517339	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	243024	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	381806	20.00	ng	0.00
76) Chrysene-d12	14.03	240	315433	20.00	ng	0.00
87) Perylene-d12	15.50	264	278728	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	372465	44.97	ng	0.00
7) Phenol-d6	6.48	99	490830	46.97	ng	0.00
23) Nitrobenzene-d5	7.41	82	278603	31.09	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	99345	42.16	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	535529	41.28	ng	0.00
79) Terphenyl-d14	12.96	244	507934	33.93	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	118026	7.133	ng	Qvalue 100
58) Fluorene	10.44	166	42833	2.995	ng	98
71) Phenanthrene	11.40	178	456851	23.406	ng	99
72) Anthracene	11.46	178	170365	8.624	ng	98
75) Fluoranthene	12.60	202	800569	39.178	ng	99
78) Pyrene	12.83	202	747107	31.420	ng	99
81) Benzo(a)anthracene	14.02	228	376549	18.677	ng	98
83) Chrysene	14.05	228	296704	15.158	ng	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	119247	7.254	ng	97
88) Benzo(b)fluoranthene	15.07	252	339157m	21.044	ng	
89) Benzo(k)fluoranthene	15.09	252	104353m	6.648	ng	
90) Benzo(a)pyrene	15.43	252	240531	16.696	ng	# 96
91) Dibenzo(a,h)anthracene	16.98	278	29418	2.359	ng	# 74
92) Benzo(a,h,i)perylene	17.43	276	117686	9.609	ng	99

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

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