

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113476.D
 Acq On : 1 Apr 2019 19:47
 Operator : JU/SJ
 Sample : K1692-05 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-032919

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:38 PM

Quant Time: Apr 02 07:26:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	135600	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	537589	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	248695	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	447787	20.00	ng	0.00
76) Chrysene-d12	13.85	240	392411	20.00	ng	0.00
87) Perylene-d12	15.26	264	275709	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	159557	19.24	ng	0.00
7) Phenol-d6	6.35	99	231039	22.02	ng	0.00
23) Nitrobenzene-d5	7.27	82	119638	13.21	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	57973	18.87	ng	0.00
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
79) Terphenyl-d14	12.79	244	279903	15.72	ng	0.00
Target Compounds						
52) Acenaphthene	9.77	154	29854	2.050	ng	Qvalue 100
71) Phenanthrene	11.24	178	291266	13.098	ng	98
72) Anthracene	11.29	178	87618	3.879	ng	97
75) Fluoranthene	12.43	202	520962	20.723	ng	100
78) Pyrene	12.65	202	484141	17.804	ng	98
81) Benzo(a)anthracene	13.84	228	238871	10.634	ng	97
83) Chrysene	13.87	228	202788	8.888	ng	98
86) Indeno(1,2,3-cd)pyrene	16.62	276	77533	3.960	ng	98
88) Benzo(b)fluoranthene	14.86	252	199937m	11.674	ng	
89) Benzo(k)fluoranthene	14.89	252	81631m	5.427	ng	
90) Benzo(a)pyrene	15.20	252	146943	9.678	ng	95
92) Benzo(g,h,i)perylene	17.03	276	83287	6.292	ng	94

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

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