

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040319\
 Data File : BF113564.D
 Acq On : 4 Apr 2019 1:18
 Operator : JU/SJ
 Sample : K2203-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-040219-A

Manual Integrations
 APPROVED

Sohil
 4/4/2019 10:58:00 AM

Quant Time: Apr 04 06:12:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	151280	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	530755	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	225911	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	404630	20.00	ng	0.00
76) Chrysene-d12	13.84	240	344791	20.00	ng	0.00
87) Perylene-d12	15.25	264	268932	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	819525	92.38	ng	0.00
7) Phenol-d6	6.33	99	1066643	97.55	ng	0.00
23) Nitrobenzene-d5	7.25	82	617794	68.18	ng	0.00
42) 2,4,6-Tribromophenol	10.51	330	223697	82.43	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1088484	77.57	ng	0.00
79) Terphenyl-d14	12.79	244	1042409	61.55	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.35	94	31821	2.548	ng	78
50) Dimethylphthalate	9.44	163	97253	5.840	ng	100
71) Phenanthrene	11.23	178	248488	12.359	ng	98
72) Anthracene	11.28	178	83260	4.070	ng	99
75) Fluoranthene	12.42	202	730574	33.909	ng	99
78) Pyrene	12.64	202	713345	27.191	ng	99
81) Benzo(a)anthracene	13.83	228	401572	20.190	ng	97
83) Chrysene	13.87	228	360842	17.897	ng	98
86) Indeno(1,2,3-cd)pyrene	16.62	276	158005	8.488	ng	93
88) Benzo(b)fluoranthene	14.86	252	425986m	25.877	ng	
89) Benzo(k)fluoranthene	14.88	252	97370m	6.591	ng	
90) Benzo(a)pyrene	15.20	252	272155	18.605	ng	99
91) Dibenzo(a,h)anthracene	16.62	278	42533	3.109	ng	# 81
92) Benzo(g,h,i)perylene	17.03	276	146123	10.465	ng	99

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

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