

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041219\
 Data File : BF113756.D
 Acq On : 12 Apr 2019 23:51
 Operator : JU/SJ
 Sample : K2376-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4-S

Manual Integrations
 APPROVED

Sohil
 4/15/2019 12:27:24 AM

Quant Time: Apr 13 04:15:25 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	6.67	152	107754	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	393055	20.00	ng	-0.01
39) Acenaphthene-d10	9.70	164	170414	20.00	ng	-0.02
64) Phenanthrene-d10	11.19	188	300834	20.00	ng	-0.02
76) Chrysene-d12	13.82	240	303843	20.00	ng	-0.02
87) Perylene-d12	15.23	264	295493	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	587463	92.97	ng	-0.01
7) Phenol-d6	6.33	99	790572	101.51	ng	-0.01
23) Nitrobenzene-d5	7.24	82	458229	68.29	ng	-0.02
42) 2,4,6-Tribromophenol	10.49	330	176629	86.28	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	767979	72.55	ng	-0.02
79) Terphenyl-d14	12.76	244	687725	46.08	ng	-0.03
Target Compounds						
49) Acenaphthylene	9.57	152	36944	2.134	ng	Qvalue 99
50) Dimethylphthalate	9.43	163	65591	5.221	ng	98
71) Phenanthrene	11.21	178	197298	13.198	ng	98
72) Anthracene	11.26	178	60929	4.006	ng	96
75) Fluoranthene	12.40	202	483121	30.160	ng	98
78) Pyrene	12.62	202	526809	22.787	ng	99
81) Benzo(a)anthracene	13.81	228	348213	19.867	ng	96
83) Chrysene	13.85	228	336628	18.946	ng	98
86) Indeno(1,2,3-cd)pyrene	16.60	276	207410	12.643	ng	93
88) Benzo(b)fluoranthene	14.84	252	486885m	26.918	ng	
89) Benzo(k)fluoranthene	14.86	252	122289m	7.533	ng	
90) Benzo(a)pyrene	15.18	252	346649	21.567	ng	100
91) Dibenzo(a,h)anthracene	16.60	278	48069	3.198	ng	# 84
92) Benzo(g,h,i)perylene	17.01	276	201787	13.153	ng	98

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

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