

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
 Data File : BF111644.D
 Acq On : 20 Dec 2018 17:47
 Operator : JU/SJ
 Sample : J6445-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EB06_12-14

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:46:46 PM

Quant Time: Dec 21 06:44:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	198606	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	737744	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	361026	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	597419	20.00	ng	0.00
76) Chrysene-d12	14.04	240	464364	20.00	ng	0.00
87) Perylene-d12	15.53	264	468522	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	769836	66.07	ng	0.00
7) Phenol-d6	6.52	99	1184428	79.88	ng	0.00
23) Nitrobenzene-d5	7.45	82	742524	58.58	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	138224	32.40	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1414787	57.12	ng	0.00
79) Terphenyl-d14	13.00	244	1148845	46.92	ng	0.00
Target Compounds						
10) Phenol	6.53	94	98277	5.874	ng	88
50) Dimethylphthalate	9.64	163	216313	8.193	ng	97
58) Fluorene	10.47	166	57345	2.423	ng	98
71) Phenanthrene	11.44	178	645157	20.363	ng	94
72) Anthracene	11.49	178	148243	4.661	ng	94
75) Fluoranthene	12.62	202	663921	20.693	ng	95
78) Pyrene	12.85	202	569449	17.365	ng	96
81) Benzo(a)anthracene	14.03	228	301210	11.366	ng	99
83) Chrysene	14.07	228	268979	10.327	ng	98
86) Indeno(1,2,3-cd)pyrene	17.00	276	131106	5.921	ng	# 83
88) Benzo(b)fluoranthene	15.09	252	356908m	13.034	ng	
89) Benzo(k)fluoranthene	15.12	252	108341m	4.156	ng	
90) Benzo(a)pyrene	15.46	252	246049m	9.891	ng	
92) Benzo(a,h,i)perylene	17.44	276	119552	5.620	ng	# 90

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

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