

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110252.D
 Acq On : 29 Oct 2018 10:37
 Operator : JU/SJ
 Sample : J5654-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW2-02

Manual Integrations
 APPROVED

Sohil
 10/30/2018 2:24:06 PM

Quant Time: Oct 29 12:40:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 12:32:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	113407	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	441494	20.00	ng	0.00
39) Acenaphthene-d10	10.01	164	190333	20.00	ng	0.00
64) Phenanthrene-d10	11.50	188	294019	20.00	ng	0.00
76) Chrysene-d12	14.16	240	209389	20.00	ng	0.00
87) Perylene-d12	15.69	264	192306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	563979	80.98	ng	0.00
7) Phenol-d6	6.59	99	907180	101.84	ng	0.00
23) Nitrobenzene-d5	7.53	82	581560	78.13	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	81069	43.00	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1038093	85.64	ng	0.00
79) Terphenyl-d14	13.09	244	736472	73.15	ng	0.00
Target Compounds						
10) Phenol	6.60	94	51800	5.440	ng	Qvalue # 63
31) Naphthalene	8.27	128	50007	2.458	ng	98
50) Dimethylphthalate	9.72	163	108675	7.735	ng	98
71) Phenanthrene	11.53	178	232863	15.136	ng	99
72) Anthracene	11.58	178	53520	3.471	ng	98
75) Fluoranthene	12.72	202	212673	13.621	ng	99
78) Pyrene	12.96	202	193641	12.900	ng	97
81) Benzo(a)anthracene	14.14	228	91561	7.374	ng	98
83) Chrysene	14.18	228	83567	7.086	ng	95
86) Indeno(1,2,3-cd)pyrene	17.26	276	31329	3.202	ng	98
88) Benzo(b)fluoranthene	15.23	252	84120m	7.575	ng	
89) Benzo(k)fluoranthene	15.26	252	33363m	3.056	ng	
90) Benzo(a)pyrene	15.62	252	61372	6.006	ng	# 96
92) Benzo(a,h,i)perylene	17.73	276	29118	3.351	ng	# 90

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

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