

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
 Data File : BF110850.D
 Acq On : 19 Nov 2018 13:58
 Operator : JU/SJ
 Sample : J5977-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-01

Manual Integrations
 APPROVED

Sohil
 11/20/2018 11:07:27 AM

Quant Time: Nov 19 15:52:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	55918	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	234162	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	110260	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	195917	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	122625	20.00	ng	-0.01
87) Perylene-d12	15.66	264	123254	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	359051	104.30	ng	0.00
7) Phenol-d6	6.57	99	457179	103.85	ng	0.00
23) Nitrobenzene-d5	7.50	82	290702	71.49	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	112440	101.21	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	560081	72.55	ng	-0.01
79) Terphenyl-d14	13.06	244	440413	67.26	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	25688	5.032	ng	Qvalue # 73
50) Dimethylphthalate	9.69	163	95465	11.603	ng	98
71) Phenanthrene	11.50	178	71502	6.812	ng	99
75) Fluoranthene	12.69	202	103541	9.839	ng	98
78) Pyrene	12.93	202	93142	9.865	ng	96
81) Benzo(a)anthracene	14.12	228	48203	6.577	ng	96
83) Chrysene	14.16	228	43179	6.184	ng	96
86) Indeno(1,2,3-cd)pyrene	17.23	276	28113	5.611	ng	99
88) Benzo(b)fluoranthene	15.20	252	60695m	8.232	ng	
89) Benzo(k)fluoranthene	15.23	252	18870m	2.590	ng	
90) Benzo(a)pyrene	15.59	252	44858	6.696	ng	96
92) Benzo(a,h,i)perylene	17.72	276	25841	4.594	ng	96

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

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