

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
 Data File : BF106982.D
 Acq On : 29 Jun 2018 19:20
 Operator : JU/SJ
 Sample : J3732-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-106-8(3-8)

Manual Integrations
 APPROVED

Sohil
 6/30/2018 10:58:09 AM

Quant Time: Jun 29 23:58:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	68128	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	247405	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	109190	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	179970	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	187630	20.00	ng	-0.02
86) Perylene-d12	15.68	264	143239	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	500921	124.50	ng	0.00
7) Phenol-d6	6.60	99	616050	122.61	ng	-0.01
23) Nitrobenzene-d5	7.52	82	383908	86.24	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	167878	132.65	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	661473	106.22	ng	-0.02
78) Terphenyl-d14	13.07	244	649992	83.91	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.70	163	88870	10.852	ng	98
70) Phenanthrene	11.51	178	56644	6.526	ng	99
71) Anthracene	11.56	178	18560	2.095	ng	99
74) Fluoranthene	12.71	202	161501	16.880	ng	93
77) Pyrene	12.94	202	147222	11.899	ng	98
80) Benzo(a)anthracene	14.13	228	109314	9.559	ng	99
82) Chrysene	14.17	228	89054	8.465	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	35466	3.972	ng	# 89
87) Benzo(b)fluoranthene	15.22	252	105299m	11.834	ng	
88) Benzo(k)fluoranthene	15.25	252	29863m	3.524	ng	
89) Benzo(a)pyrene	15.62	252	67712m	8.560	ng	
91) Benzo(a,h,i)perylene	17.74	276	30486	4.653	ng	# 90

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

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