

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082817\
 Data File : BF098064.D
 Acq On : 28 Aug 2017 13:08
 Operator : SJ/JU
 Sample : I4961-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-COMP

Quant Time: Aug 28 14:02:49 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	107005	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	426683	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	187970	20.00	ng	-0.01
63) Phenanthrene-d10	11.26	188	351943	20.00	ng	0.00
75) Chrysene-d12	13.90	240	230493	20.00	ng	0.00
86) Perylene-d12	15.33	264	232925	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	894799	127.20	ng	0.00
7) Phenol-d6	6.38	99	1250222	130.80	ng	0.00
23) Nitrobenzene-d5	7.31	82	779988	99.31	ng	-0.01
41) 2,4,6-Tribromophenol	10.57	330	209135	128.23	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1106837	86.51	ng	0.00
78) Terphenyl-d14	12.85	244	815751	73.80	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.49	163	124845	9.086	ng	99
70) Phenanthrene	11.28	178	101895	5.433	ng	99
71) Anthracene	11.33	178	57928	3.045	ng	98
74) Fluoranthene	12.47	202	818340	41.806	ng	98
77) Pyrene	12.72	202	1852784	98.282	ng	99
80) Benzo(a)anthracene	13.89	228	813743	58.905	ng	97
82) Chrysene	13.93	228	776456	56.502	ng	98
83) Bis(2-ethylhexyl)phthalate	13.89	149	29475	3.680	ng	# 93
85) Indeno(1,2,3-cd)pyrene	16.72	276	266479	26.360	ng	# 90
87) Benzo(b)fluoranthene	14.93	252	706328m	48.108	ng	
88) Benzo(k)fluoranthene	14.94	252	270959m	19.644	ng	
89) Benzo(a)pyrene	15.27	252	739451	56.891	ng	99
90) Dibenzo(a,h)anthracene	16.73	278	89580	8.466	ng	96
91) Benzo(g,h,i)perylene	17.14	276	273398	24.762	ng	93

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

