

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090117\
 Data File : BF098218.D
 Acq On : 1 Sep 2017 11:24
 Operator : SJ/JU
 Sample : I5027-01 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-02-082917

Manual Integrations
 APPROVED

mohammad
 9/5/2017 5:21:44 PM

Quant Time: Sep 01 13:13:57 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 01 12:59:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	123174	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	493026	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	214951	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	364022	20.00	ng	0.00
75) Chrysene-d12	13.87	240	255670	20.00	ng	0.00
86) Perylene-d12	15.28	264	256592	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	185360	22.89	ng	0.00
7) Phenol-d6	6.36	99	249801	22.70	ng	-0.01
23) Nitrobenzene-d5	7.28	82	149300	16.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	37579	20.15	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	249433	17.05	ng	0.00
78) Terphenyl-d14	12.82	244	154137	12.57	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	49386	3.143	ng	99
70) Phenanthrene	11.26	178	122046	6.292	ng	98
74) Fluoranthene	12.44	202	245494	12.125	ng	98
77) Pyrene	12.67	202	203118	9.714	ng	98
80) Benzo(a)anthracene	13.86	228	78898	5.149	ng	99
82) Chrysene	13.89	228	86732	5.690	ng	95
83) Bis(2-ethylhexyl)phthalate	13.85	149	26739	3.010	ng	98
85) Indeno(1,2,3-cd)pyrene	16.64	276	51583	4.600	ng	# 90
87) Benzo(b)fluoranthene	14.87	252	114814m	7.099	ng	
88) Benzo(k)fluoranthene	14.90	252	42429m	2.792	ng	
89) Benzo(a)pyrene	15.22	252	82238	5.744	ng	98
91) Benzo(a,h,i)perylene	17.06	276	51782	4.257	ng	# 92

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

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