

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071118\
 Data File : BF107310.D
 Acq On : 11 Jul 2018 19:46
 Operator : JU/SJ
 Sample : J3914-07 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NEMH

Manual Integrations
 APPROVED

Sohil
 7/12/2018 2:19:42 PM

Quant Time: Jul 12 02:26:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 11 16:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	66114	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	246448	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	110328	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	195586	20.00	ng	0.00
75) Chrysene-d12	14.14	240	173815	20.00	ng	0.00
86) Perylene-d12	15.68	264	119788	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	175844	45.04	ng	0.00
7) Phenol-d6	6.58	99	222745	45.68	ng	0.00
23) Nitrobenzene-d5	7.50	82	127946	28.85	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	51228	40.06	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	243938	27.43	ng	0.00
78) Terphenyl-d14	13.07	244	244044	34.01	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	44426	5.369	ng	99
70) Phenanthrene	11.50	178	90527	9.596	ng	98
71) Anthracene	11.55	178	21649	2.248	ng	96
74) Fluoranthene	12.70	202	177652	17.086	ng	96
77) Pyrene	12.94	202	163030	14.224	ng	99
80) Benzo(a)anthracene	14.13	228	85322m	8.054	ng	
82) Chrysene	14.17	228	88810m	9.112	ng	
85) Indeno(1,2,3-cd)pyrene	17.28	276	38013	4.596	ng	# 86
87) Benzo(b)fluoranthene	15.22	252	92849m	12.478	ng	
88) Benzo(k)fluoranthene	15.25	252	25182m	3.554	ng	
89) Benzo(a)pyrene	15.62	252	57605	8.708	ng	98
91) Benzo(a,h,i)perylene	17.77	276	38167	6.965	ng	# 87

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

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