

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\
 Data File : BF088905.D
 Acq On : 11 Jul 2016 17:58
 Operator : UM/SJ
 Sample : H3915-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
Client Sampled :
 P7-B2-0-5-C

Manual Integrations
 APPROVED
 sohil
 7/12/2016 7:36:40 PM

Quant Time: Jul 12 02:15:15 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 11 18:13:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	6.71	152	63154	20.00	ng	0.00	
21) Naphthalene-d8	8.00	136	268294	20.00	ng	0.00	
38) Acenaphthene-d10	9.75	164	109096	20.00	ng	0.00	
63) Phenanthrene-d10	11.22	188	186326	20.00	ng	0.00	
75) Chrysene-d12	13.85	240	140621	20.00	ng	0.00	
86) Perylene-d12	15.22	264	131157	20.00	ng	0.01	
System Monitoring Compounds							
5) 2-Fluorophenol	5.30	112	371875	88.25	ng	0.01	
7) Phenol-d6	6.35	99	478837	87.94	ng	0.00	
23) Nitrobenzene-d5	7.28	82	358123	69.95	ng	0.00	
41) 2,4,6-Tribromophenol	10.52	330	99869	98.58	ng	-0.01	
44) 2-Fluorobiphenyl	9.07	172	525743	76.12	ng	0.00	
78) Terphenyl-d14	12.81	244	428709	67.90	ng	0.01	
Target Compounds							
49) Dimethylphthalate	9.47	163	82926	10.67	ng		98
70) Phenanthrene	11.24	178	87736	8.61	ng		97
74) Fluoranthene	12.42	202	223019	21.60	ng		97
77) Pyrene	12.65	202	305076	25.59	ng		99
80) Benzo(a)anthracene	13.84	228	204760	22.61	ng		100
82) Chrysene	13.87	228	199327	22.25	ng		99
85) Indeno(1,2,3-cd)pyrene	16.50	276	59849	8.68	ng	#	100
87) Benzo(b)fluoranthene	14.85	252	182454m	22.18	ng		
88) Benzo(k)fluoranthene	14.86	252	32500m	4.54	ng		
89) Benzo(a)pyrene	15.17	252	119279m	17.12	ng		
90) Dibenzo(a,h)anthracene	16.51	278	18757	3.22	ng		98
91) Benzo(a,h,i)perylene	16.88	276	61278	10.18	ng		97

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

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