

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070117\
 Data File : BF096389.D
 Acq On : 1 Jul 2017 20:04
 Operator : SJ/MA
 Sample : I3908-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-61-1.5-3

Manual Integrations
 APPROVED

mohammad
 7/3/2017 4:02:56 PM

Quant Time: Jul 03 02:42:14 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 01:43:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	149131	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	581233	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	201566	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	279234	20.00	ng	0.00
75) Chrysene-d12	13.87	240	260504	20.00	ng	0.00
86) Perylene-d12	15.29	264	233904	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	898810	88.96	ng	0.00
7) Phenol-d6	6.37	99	1000535	80.55	ng	0.00
23) Nitrobenzene-d5	7.29	82	657442	67.97	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	134600	85.49	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	857118	72.70	ng	0.00
78) Terphenyl-d14	12.83	244	502031	41.18	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	8.03	128	266038	9.695	ng	99
37) 2-Methylnaphthalene	8.73	142	153893	8.811	ng	97
45) 1,1'-Biphenyl	9.19	154	39483	2.286	ng	92
48) Acenaphthylene	9.62	152	82766	4.058	ng	98
49) Dimethylphthalate	9.48	163	256228	17.028	ng	99
51) Acenaphthene	9.80	154	60845	4.840	ng	99
54) Dibenzofuran	9.97	168	113124	6.816	ng	94
57) Fluorene	10.31	166	54744	4.670	ng	# 91
70) Phenanthrene	11.27	178	446813	29.601	ng	99
71) Anthracene	11.32	178	145310	9.447	ng	99
72) Carbazole	11.47	167	61431	3.971	ng	97
73) Di-n-butylphthalate	11.82	149	39987	2.134	ng	97
74) Fluoranthene	12.46	202	583889	39.453	ng	98
77) Pyrene	12.68	202	536569	25.661	ng	99
80) Benzo(a)anthracene	13.87	228	297282	19.707	ng	94
82) Chrysene	13.90	228	357867	22.653	ng	97
83) Bis(2-ethylhexyl)phthalate	13.87	149	103353	8.747	ng	# 97
85) Indeno(1,2,3-cd)pyrene	16.66	276	132959	10.663	ng	98
87) Benzo(b)fluoranthene	14.89	252	526091m	35.895	ng	
88) Benzo(k)fluoranthene	14.92	252	118287m	9.022	ng	
89) Benzo(a)pyrene	15.23	252	219782	16.796	ng	96
90) Dibenz(a,h)anthracene	16.67	278	38259	3.717	ng	# 82
91) Benzo(g,h,i)perylene	17.07	276	117332	11.000	ng	97

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

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