

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079732.D
 Acq On : 5 Jun 2015 7:37
 Operator : TP/IZ
 Sample : G2450-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEC-3.4-SP

Quant Time: Jun 06 04:33:36 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 06 03:31:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	64603	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	253507	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	141396	20.00	ng	0.00
63) Phenanthrene-d10	12.09	188	273697	20.00	ng	0.01
75) Chrysene-d12	15.12	240	292405	20.00	ng	0.14
86) Perylene-d12	17.10	264	294133	20.00	ng	0.18

System Monitoring Compounds						
5) 2-Fluorophenol	6.14	112	444218	117.90	ng	0.01
7) Phenol-d6	7.11	99	625270	127.49	ng	0.01
23) Nitrobenzene-d5	8.07	82	319439	76.72	ng	0.00
41) 2,4,6-Tribromophenol	11.38	330	159402	125.54	ng	0.01
44) 2-Fluorobiphenyl	9.87	172	732421	74.99	ng	0.00
78) Terphenyl-d14	13.81	244	946933	76.30	ng	0.07

Target Compounds				Qvalue		
31) Naphthalene	8.82	128	97049	6.95	ng	99
37) 2-Methylnaphthalene	9.52	142	32830	3.48	ng	96
48) Acenaphthylene	10.44	152	188530	11.92	ng	99
49) Dimethylphthalate	10.26	163	70790	6.50	ng	98
51) Acenaphthene	10.62	154	22595	2.43	ng	92
54) Dibenzofuran	10.79	168	37560	2.81	ng	98
57) Fluorene	11.13	166	36655	3.22	ng	# 96
70) Phenanthrene	12.13	178	491769	30.89	ng	99
71) Anthracene	12.17	178	339640	21.56	ng	99
72) Carbazole	12.32	167	136915	9.32	ng	97
74) Fluoranthene	13.41	202	1820301	105.40	ng	99
77) Pyrene	13.67	202	1661043	91.42	ng	99
80) Benzo(a)anthracene	15.10	228	972330	54.39	ng	97
82) Chrysene	15.15	228	973886	58.22	ng	98
85) Indeno(1,2,3-cd)pyrene	19.12	276	446549	23.65	ng	95
87) Benzo(b)fluoranthene	16.51	252	1244612m	65.60	ng	
88) Benzo(k)fluoranthene	16.55	252	428871m	23.69	ng	
89) Benzo(a)pyrene	17.02	252	695945	40.15	ng	97
90) Dibenzo(a,h)anthracene	19.13	278	123127	7.69	ng	# 87
91) Benzo(g,h,i)perylene	19.74	276	401491	24.52	ng	99

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

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