

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079734.D
 Acq On : 5 Jun 2015 10:01
 Operator : TP/IZ
 Sample : G2456-01DL 2X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILL-02-ALLOCCODL

Quant Time: Jun 08 03:14:33 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 19:33:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	56105	20.00	ng	0.01
21) Naphthalene-d8	8.81	136	224426	20.00	ng	0.01
38) Acenaphthene-d10	10.58	164	128896	20.00	ng	0.00
63) Phenanthrene-d10	12.11	188	261655	20.00	ng	0.01
75) Chrysene-d12	15.27	240	272937	20.00	ng	0.10
86) Perylene-d12	17.33	264	237279	20.00	ng	0.16

System Monitoring Compounds

5) 2-Fluorophenol	6.12	112	89158	27.25	ng	0.00
7) Phenol-d6	7.11	99	193208	45.36	ng	0.01
23) Nitrobenzene-d5	8.07	82	94879	25.74	ng	0.00
41) 2,4,6-Tribromophenol	11.38	330	3750	3.24	ng	0.01
44) 2-Fluorobiphenyl	9.87	172	258272	29.01	ng	0.00
78) Terphenyl-d14	13.89	244	355910	30.72	ng	0.06

Target Compounds

						Qvalue
31) Naphthalene	8.82	128	141094	11.41	ng	99
37) 2-Methylnaphthalene	9.52	142	48171	5.77	ng	98
49) Dimethylphthalate	10.26	163	23888	2.41	ng	98
51) Acenaphthene	10.62	154	51670	6.09	ng	98
54) Dibenzofuran	10.79	168	85451	7.02	ng	98
57) Fluorene	11.13	166	95531	9.19	ng	95
70) Phenanthrene	12.14	178	656572	43.14	ng	99
71) Anthracene	12.19	178	258254	17.15	ng	100
72) Carbazole	12.34	167	89738	6.39	ng	98
74) Fluoranthene	13.46	202	547897	33.19	ng	98
77) Pyrene	13.74	202	452713	26.69	ng	99
79) Butylbenzylphthalate	14.47	149	14533	2.22	ng	# 96
80) Benzo(a)anthracene	15.25	228	233729	14.01	ng	99
82) Chrysene	15.30	228	205206	13.14	ng	95
83) Bis(2-ethylhexyl)phthalate	15.21	149	23748	2.31	ng	# 89
85) Indeno(1,2,3-cd)pyrene	19.37	276	78163	4.44	ng	95
87) Benzo(b)fluoranthene	16.73	252	219336m	14.33	ng	
88) Benzo(k)fluoranthene	16.77	252	70962m	4.86	ng	
89) Benzo(a)pyrene	17.23	252	144144	10.31	ng	# 90
91) Benzo(g,h,i)perylene	19.99	276	75567	5.72	ng	94

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

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