

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079729.D
 Acq On : 5 Jun 2015 6:10
 Operator : TP/IZ
 Sample : G2456-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILL-02-ALLOCCO

Quant Time: Jun 06 04:21:26 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	62440	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	244488	20.00	ng	0.00
38) Acenaphthene-d10	10.58	164	121191	20.00	ng	0.00
63) Phenanthrene-d10	12.10	188	255417	20.00	ng	0.02
75) Chrysene-d12	15.19	240	280435m	20.00	ng	0.21
86) Perylene-d12	17.21	264	258459m	20.00	ng	0.30

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	6.12	112	188383	51.73	ng	0.00
7) Phenol-d6	7.10	99	374205	78.94	ng	0.00
23) Nitrobenzene-d5	8.06	82	193720	48.24	ng	-0.01
41) 2,4,6-Tribromophenol	11.37	330	7335	6.74	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	541805	64.73	ng	0.00
78) Terphenyl-d14	13.84	244	669398	56.24	ng	0.10

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.82	128	303883	22.56	ng	99
37) 2-Methylnaphthalene	9.52	142	103351	11.37	ng	98
45) 1,1'-Biphenyl	9.98	154	26264	2.71	ng	91
49) Dimethylphthalate	10.26	163	43827	4.69	ng	# 98
51) Acenaphthene	10.60	154	98371	12.33	ng	97
54) Dibenzofuran	10.79	168	155398	13.59	ng	95
57) Fluorene	11.13	166	208528	21.34	ng	99
70) Phenanthrene	12.12	178	1231636	82.90	ng	99
71) Anthracene	12.18	178	497621	33.85	ng	100
72) Carbazole	12.33	167	173955	12.69	ng	99
74) Fluoranthene	13.43	202	1062155	65.90	ng	100
77) Pyrene	13.69	202	886681	50.88	ng	99
80) Benzo(a)anthracene	15.18	228	460496m	26.86	ng	
82) Chrysene	15.22	228	432236m	26.94	ng	
83) Bis(2-ethylhexyl)phthalate	15.13	149	43501m	4.13	ng	
85) Indeno(1,2,3-cd)pyrene	19.26	276	148894m	8.22	ng	
87) Benzo(b)fluoranthene	16.62	252	409119m	24.54	ng	
88) Benzo(k)fluoranthene	16.64	252	180352m	11.34	ng	
89) Benzo(a)pyrene	17.13	252	310942m	20.41	ng	
90) Dibenzo(a,h)anthracene	19.27	278	46690m	3.32	ng	
91) Benzo(g,h,i)perylene	19.87	276	145844m	10.14	ng	

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

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