

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078577.D
 Acq On : 16 Apr 2015 20:46
 Operator : TP/IZ
 Sample : G1855-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LS-SB16A(6-12)

Quant Time: Apr 17 02:26:35 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 00:52:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	20626	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	90589	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	46617	20.00	ng	0.00
63) Phenanthrene-d10	12.24	188	93232	20.00	ng	0.00
75) Chrysene-d12	15.50	240	107079	20.00	ng	0.01
86) Perylene-d12	17.22	264	116847	20.00	ng	0.03

System Monitoring Compounds

5) 2-Fluorophenol	4.94	112	125524m	100.34	ng	0.03
7) Phenol-d6	6.31	99	165175	105.36	ng	0.01
23) Nitrobenzene-d5	7.39	82	99907	66.85	ng	0.00
41) 2,4,6-Tribromophenol	11.41	330	41085	106.27	ng	0.01
44) 2-Fluorobiphenyl	9.62	172	206141	63.21	ng	0.00
78) Terphenyl-d14	14.24	244	231633	48.88	ng	0.01

Target Compounds

						Qvalue
31) Naphthalene	8.30	128	22645	4.80	ng	98
37) 2-Methylnaphthalene	9.15	142	10192	3.18	ng	# 90
48) Acenaphthylene	10.25	152	21227	4.38	ng	99
49) Dimethylphthalate	10.14	163	14937	4.26	ng	# 99
54) Dibenzofuran	10.67	168	11228	2.70	ng	96
57) Fluorene	11.10	166	7775	2.30	ng	# 94
70) Phenanthrene	12.27	178	149568	27.73	ng	99
71) Anthracene	12.33	178	46385	8.73	ng	99
72) Carbazole	12.55	167	12804	2.57	ng	# 92
74) Fluoranthene	13.74	202	290560	51.09	ng	92
77) Pyrene	14.00	202	266408	39.63	ng	# 93
80) Benzo(a)anthracene	15.49	228	173262	27.20	ng	95
82) Chrysene	15.53	228	145520	24.31	ng	97
83) Bis(2-ethylhexyl)phthalate	15.60	149	512956	127.66	ng	98
85) Indeno(1,2,3-cd)pyrene	18.45	276	129411m	19.05	ng	
87) Benzo(b)fluoranthene	16.79	252	231149m	32.01	ng	
88) Benzo(k)fluoranthene	16.81	252	86450m	13.47	ng	
89) Benzo(a)pyrene	17.17	252	175047	27.77	ng	97
90) Dibenzo(a,h)anthracene	18.46	278	34160	5.41	ng	# 85
91) Benzo(g,h,i)perylene	18.77	276	117832	18.63	ng	96

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

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