

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041315\
 Data File : BF078498.D
 Acq On : 14 Apr 2015 11:19
 Operator : TP/IZ
 Sample : G1850-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-ODA

Manual Integrations
 APPROVED

apatel
 4/14/2015 5:27:08 PM

Quant Time: Apr 14 12:09:01 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 14 11:28:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	33380	20.00	ng	0.00
21) Naphthalene-d8	8.41	136	137821	20.00	ng	0.00
38) Acenaphthene-d10	10.56	164	69057	20.00	ng	-0.01
63) Phenanthrene-d10	12.39	188	132466	20.00	ng	0.00
75) Chrysene-d12	15.63	240	144811	20.00	ng	-0.01
86) Perylene-d12	17.34	264	149581	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.08	112	201685	99.62	ng	0.01
7) Phenol-d6	6.43	99	271500	107.01	ng	0.00
23) Nitrobenzene-d5	7.53	82	153343	67.44	ng	0.00
41) 2,4,6-Tribromophenol	11.54	330	65387	114.17	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	283417	58.66	ng	0.00
78) Terphenyl-d14	14.38	244	324862	50.69	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.44	94	6564	2.16	ng	# 61
31) Naphthalene	8.43	128	150027	20.91	ng	98
37) 2-Methylnaphthalene	9.29	142	26386	5.42	ng	99
48) Acenaphthylene	10.39	152	14846	2.07	ng	98
49) Dimethylphthalate	10.26	163	15842	3.05	ng	100
51) Acenaphthene	10.60	154	67753	16.77	ng	100
54) Dibenzofuran	10.82	168	23403	3.79	ng	97
57) Fluorene	11.23	166	36335	7.26	ng	98
70) Phenanthrene	12.41	178	198238	25.87	ng	99
71) Anthracene	12.48	178	89062	11.80	ng	99
72) Carbazole	12.70	167	14557	2.06	ng	94
74) Fluoranthene	13.87	202	238089	29.46	ng	96
77) Pyrene	14.15	202	325489	35.80	ng	97
80) Benzo(a)anthracene	15.62	228	118111m	13.71	ng	
82) Chrysene	15.67	228	79266	9.79	ng	97
85) Indeno(1,2,3-cd)pyrene	18.56	276	47005m	5.12	ng	
87) Benzo(b)fluoranthene	16.90	252	103463m	11.19	ng	
88) Benzo(k)fluoranthene	16.93	252	20431m	2.49	ng	
89) Benzo(a)pyrene	17.27	252	101755	12.61	ng	# 92
91) Benzo(g,h,i)perylene	18.91	276	54347	6.71	ng	96

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

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