

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
 Data File : BF086691.D
 Acq On : 4 May 2016 22:05
 Operator : UM/SJ
 Sample : H2715-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-3-CS-9(0-12FTBGS)

Manual Integrations
 APPROVED

sohil
 5/5/2016 3:20:06 PM

Quant Time: May 05 04:52:08 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 15:12:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	157173	40.00	ng	0.00
21) Naphthalene-d8	8.03	136	588617	40.00	ng	0.00
38) Acenaphthene-d10	9.78	164	249675	40.00	ng	0.00
63) Phenanthrene-d10	11.25	188	360943m	40.00	ng	0.00
75) Chrysene-d12	13.89	240	329617	40.00	ng	0.01
86) Perylene-d12	15.28	264	252221	40.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1211223	127.58	ng	0.01
7) Phenol-d6	6.38	99	1565316	122.40	ng	0.01
23) Nitrobenzene-d5	7.31	82	1196255	104.44	ng	0.00
41) 2,4,6-Tribromophenol	10.57	330	309367	126.67	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1457061	95.66	ng	0.00
78) Terphenyl-d14	12.84	244	1023223	62.74	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.04	128	69639	2.32	ng	99
49) Dimethylphthalate	9.49	163	158716	8.53	ng	98
57) Fluorene	10.32	166	43569	2.81	ng	96
70) Phenanthrene	11.28	178	537715	27.73	ng	99
71) Anthracene	11.33	178	91115	4.49	ng	99
72) Carbazole	11.49	167	51875	2.94	ng	99
74) Fluoranthene	12.46	202	592042	30.47	ng	98
77) Pyrene	12.69	202	560097	21.31	ng	99
80) Benzo(a)anthracene	13.88	228	220204	11.81	ng	95
82) Chrysene	13.92	228	240132	12.07	ng	98
85) Indeno(1,2,3-cd)pyrene	16.60	276	79441	4.66	ng	# 92
87) Benzo(b)fluoranthene	14.89	252	212106m	14.00	ng	
88) Benzo(k)fluoranthene	14.90	252	115139m	8.06	ng	
89) Benzo(a)pyrene	15.22	252	156908	11.24	ng	# 92
91) Benzo(a,h,i)perylene	16.99	276	75082	5.33	ng	96

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

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