

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080315\
 Data File : BF080834.D
 Acq On : 4 Aug 2015 10:15
 Operator : TP
 Sample : G3145-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-003-A

Manual Integrations
 APPROVED

MMDadoda
 8/5/2015 2:29:28 PM

Quant Time: Aug 04 11:33:14 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 01:58:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	175499	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	601695	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	258374	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	392399	20.00	ng	0.00
75) Chrysene-d12	13.97	240	343884	20.00	ng	-0.01
86) Perylene-d12	15.39	264	375715	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1515004	143.92	ng	0.01
7) Phenol-d6	6.46	99	1959161	138.36	ng	0.00
23) Nitrobenzene-d5	7.40	82	1381701	113.32	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	335890	135.66	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	1790564	104.75	ng	0.00
78) Terphenyl-d14	12.93	244	1043255	60.46	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.58	163	208703	12.09	ng	98
70) Phenanthrene	11.37	178	153706	7.53	ng	98
74) Fluoranthene	12.56	202	231851	11.86	ng	92
77) Pyrene	12.78	202	195405	7.67	ng	96
80) Benzo(a)anthracene	13.96	228	142780	7.27	ng	97
82) Chrysene	14.00	228	100405	5.26	ng	97
85) Indeno(1,2,3-cd)pyrene	16.75	276	57646	3.29	ng	95
87) Benzo(b)fluoranthene	14.99	252	157600m	7.17	ng	
88) Benzo(k)fluoranthene	15.01	252	49192m	2.57	ng	
89) Benzo(a)pyrene	15.33	252	103852	5.54	ng	97
91) Benzo(g,h,i)perylene	17.16	276	54897	2.90	ng	# 88

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

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