

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021016\
 Data File : BF084992.D
 Acq On : 10 Feb 2016 14:48
 Operator : SJ/IZ
 Sample : H1392-05 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STOCKPILE-7E-(0-2)B

Manual Integrations
 APPROVED

UMANGI
 2/11/2016 10:46:40 AM

Quant Time: Feb 10 16:01:53 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 10 13:28:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	148380	20.00	ng	-0.01
21) Naphthalene-d8	7.91	136	581414	20.00	ng	-0.01
38) Acenaphthene-d10	9.66	164	237662	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	402261	20.00	ng	0.00
75) Chrysene-d12	13.77	240	309555	20.00	ng	0.00
86) Perylene-d12	15.13	264	286128	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	256730	26.95	ng	0.00
7) Phenol-d6	6.27	99	308274	26.10	ng	-0.01
23) Nitrobenzene-d5	7.19	82	196226	19.01	ng	-0.01
41) 2,4,6-Tribromophenol	10.44	330	58442	23.57	ng	-0.01
44) 2-Fluorobiphenyl	8.99	172	303621	16.17	ng	0.00
78) Terphenyl-d14	12.72	244	207597	15.20	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.39	163	65054	3.59	ng	# 90
70) Phenanthrene	11.17	178	81180	3.64	ng	# 93
74) Fluoranthene	12.34	202	133933	5.93	ng	98
77) Pyrene	12.57	202	151712	6.40	ng	98
80) Benzo(a)anthracene	13.76	228	71515	3.72	ng	94
82) Chrysene	13.79	228	62969m	3.38	ng	
85) Indeno(1,2,3-cd)pyrene	16.38	276	33961	2.32	ng	97
87) Benzo(b)fluoranthene	14.75	252	88730m	4.62	ng	
89) Benzo(a)pyrene	15.07	252	56362	3.44	ng	# 88
91) Benzo(g,h,i)perylene	16.75	276	32663	2.35	ng	# 92

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

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