

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
 Data File : BF087799.D
 Acq On : 7 Jun 2016 21:50
 Operator : UM/SJ
 Sample : H3379-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-1

Manual Integrations
 APPROVED

sohil
 6/8/2016 7:13:59 PM

Quant Time: Jun 08 12:04:34 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	93729	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	423727	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	196048	20.00	ng	-0.01
63) Phenanthrene-d10	11.35	188	331907	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	237352	20.00	ng	-0.01
86) Perylene-d12	15.39	264	216049	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	511851	93.36	ng	0.01
7) Phenol-d6	6.47	99	558091	83.99	ng	-0.01
23) Nitrobenzene-d5	7.40	82	413429	56.97	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	172747	83.45	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	683363	52.65	ng	-0.01
78) Terphenyl-d14	12.94	244	605385	58.04	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.59	163	102177	7.19	ng	99
70) Phenanthrene	11.37	178	122220	7.02	ng	99
74) Fluoranthene	12.56	202	200943	10.93	ng	95
77) Pyrene	12.79	202	185638	10.54	ng	97
80) Benzo(a)anthracene	13.97	228	86994	6.43	ng	99
82) Chrysene	14.00	228	59873	4.71	ng	95
85) Indeno(1,2,3-cd)pyrene	16.77	276	39937	3.38	ng	# 100
87) Benzo(b)fluoranthene	14.98	252	92160m	6.12	ng	
89) Benzo(a)pyrene	15.33	252	61197	5.02	ng	94
91) Benzo(g,h,i)perylene	17.18	276	41997	3.61	ng	95

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

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