

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\
 Data File : BF087856.D
 Acq On : 9 Jun 2016 17:00
 Operator : UM/SJ
 Sample : H3399-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-25

Quant Time: Jun 09 17:49:22 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	6.82	152	142179	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	503092	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	224270	20.00	ng	-0.03
63) Phenanthrene-d10	11.34	188	358190	20.00	ng	-0.02
75) Chrysene-d12	13.95	240	275703	20.00	ng	-0.03
86) Perylene-d12	15.36	264	257053	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	368243	44.28	ng	-0.02
7) Phenol-d6	6.44	99	444448	44.10	ng	-0.03
23) Nitrobenzene-d5	7.38	82	306164	35.54	ng	-0.03
41) 2,4,6-Tribromophenol	10.64	330	97761	41.28	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	484824	30.40	ng	-0.03
78) Terphenyl-d14	12.91	244	333477	27.52	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	8.12	128	57930	2.33	ng	98
49) Dimethylphthalate	9.58	163	36819	2.27	ng	# 98
51) Acenaphthene	9.88	154	39575	2.75	ng	98
70) Phenanthrene	11.36	178	434800	23.13	ng	98
71) Anthracene	11.40	178	124062	6.54	ng	98
72) Carbazole	11.55	167	51579	2.91	ng	98
74) Fluoranthene	12.54	202	504763	25.45	ng	99
77) Pyrene	12.77	202	459021	22.44	ng	99
80) Benzo(a)anthracene	13.94	228	273470	17.41	ng	98
82) Chrysene	13.98	228	187263	12.67	ng	96
83) Bis(2-ethylhexyl)phthalate	13.94	149	35536	3.23	ng	98
85) Indeno(1,2,3-cd)pyrene	16.72	276	104438	7.61	ng	# 100
87) Benzo(b)fluoranthene	14.96	252	302004m	16.86	ng	
88) Benzo(k)fluoranthene	14.98	252	62184m	4.60	ng	
89) Benzo(a)pyrene	15.30	252	190013	13.11	ng	99
91) Benzo(g,h,i)perylene	17.13	276	93854	6.78	ng	96

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

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