

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\
 Data File : BF087854.D
 Acq On : 9 Jun 2016 16:02
 Operator : UM/SJ
 Sample : H3399-16 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-35

Quant Time: Jun 09 16:57:41 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	127815	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	468666	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	209756	20.00	ng	-0.03
63) Phenanthrene-d10	11.32	188	335196	20.00	ng	-0.03
75) Chrysene-d12	13.95	240	244398	20.00	ng	-0.03
86) Perylene-d12	15.36	264	247366	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	217891	29.14	ng	-0.02
7) Phenol-d6	6.44	99	289116	31.91	ng	-0.03
23) Nitrobenzene-d5	7.38	82	151080	18.82	ng	-0.03
41) 2,4,6-Tribromophenol	10.64	330	59106	26.69	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	304017	18.43	ng	-0.03
78) Terphenyl-d14	12.91	244	175976	16.38	ng	-0.03
Target Compounds						Qvalue
70) Phenanthrene	11.35	178	89277	5.08	ng	98
74) Fluoranthene	12.54	202	122257	6.59	ng	94
77) Pyrene	12.76	202	102193	5.63	ng	97
80) Benzo(a)anthracene	13.94	228	69390	4.98	ng	97
82) Chrysene	13.98	228	47401	3.62	ng	96
85) Indeno(1,2,3-cd)pyrene	16.71	276	30650	2.52	ng	# 100
87) Benzo(b)fluoranthene	14.96	252	73026m	4.24	ng	
88) Benzo(k)fluoranthene	14.97	252	28797m	2.21	ng	
89) Benzo(a)pyrene	15.25	252	39265	2.81	ng	96
91) Benzo(g,h,i)perylene	17.12	276	29359	2.20	ng	96

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

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