

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051916\
 Data File : BF087190.D
 Acq On : 19 May 2016 16:42
 Operator : UM/SJ
 Sample : H3119-01DL 100X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-CLARK-ST-ESMI-7DL

Manual Integrations
 APPROVED

umangi
 5/20/2016 7:13:16 PM

Quant Time: May 19 18:37:25 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF051716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 19 18:14:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	126956	20.00	ng	-0.01
21) Naphthalene-d8	7.85	136	547155	20.00	ng	-0.01
38) Acenaphthene-d10	9.60	164	271809	20.00	ng	0.00
63) Phenanthrene-d10	11.07	188	502539	20.00	ng	-0.01
75) Chrysene-d12	13.70	240	403311	20.00	ng	-0.01
86) Perylene-d12	15.05	264	293752	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	0.00	112	0	0.00	ng	
7) Phenol-d6	6.27	99	7096	0.70	ng	0.03
23) Nitrobenzene-d5	7.16	82	8010	0.73	ng	0.01
41) 2,4,6-Tribromophenol	10.40	330	214	0.07	ng	0.00
44) 2-Fluorobiphenyl	8.94	172	18281	1.11	ng	0.00
78) Terphenyl-d14	12.66	244	15996	0.90	ng	0.00
Target Compounds						
31) Naphthalene	7.87	128	1618361	60.53	ng	98
37) 2-Methylnaphthalene	8.57	142	241975	14.15	ng	99
45) 1,1'-Biphenyl	9.03	154	89801	3.98	ng	96
51) Acenaphthene	9.63	154	341478	20.69	ng	99
57) Fluorene	10.15	166	182790	10.66	ng	99
70) Phenanthrene	11.11	178	932932	34.23	ng	99
71) Anthracene	11.15	178	371980	13.49	ng	99
74) Fluoranthene	12.29	202	400092	14.65	ng	92
77) Pyrene	12.51	202	535041	18.36	ng	98
80) Benzo(a)anthracene	13.69	228	129264m	5.54	ng	
82) Chrysene	13.73	228	92316	4.09	ng	98
87) Benzo(b)fluoranthene	14.70	252	66950m	3.44	ng	
89) Benzo(a)pyrene	15.01	252	78416	4.81	ng	98
91) Benzo(g,h,i)perylene	16.65	276	36693	2.65	ng	94

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

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