

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072916\
 Data File : BF089381.D
 Acq On : 29 Jul 2016 00:10
 Operator : UM/SJ
 Sample : H4205-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-38-0.5-1.5

Quant Time: Jul 29 03:09:05 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 16:02:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	46414	20.00	ng	-0.02
21) Naphthalene-d8	7.80	136	192579	20.00	ng	-0.02
38) Acenaphthene-d10	9.55	164	89577	20.00	ng	-0.02
63) Phenanthrene-d10	11.03	188	144982	20.00	ng	-0.02
75) Chrysene-d12	13.65	240	95582	20.00	ng	-0.02
86) Perylene-d12	14.98	264	100333	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.10	112	282711	97.26	ng	-0.01
7) Phenol-d6	6.18	99	375180	97.66	ng	-0.02
23) Nitrobenzene-d5	7.10	82	244767	75.02	ng	-0.02
41) 2,4,6-Tribromophenol	10.34	330	67082	90.42	ng	-0.02
44) 2-Fluorobiphenyl	8.89	172	441268	72.29	ng	-0.02
78) Terphenyl-d14	12.60	244	261004	63.91	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.29	163	118276	17.35	ng	100
70) Phenanthrene	11.05	178	43150	5.77	ng	99
74) Fluoranthene	12.23	202	78141	9.93	ng	99
77) Pyrene	12.44	202	68894m	9.51	ng	
80) Benzo(a)anthracene	13.63	228	31573	5.85	ng	96
82) Chrysene	13.67	228	25148	4.38	ng	98
85) Indeno(1,2,3-cd)pyrene	16.17	276	17891	4.38	ng	95
87) Benzo(b)fluoranthene	14.63	252	33296m	5.34	ng	
88) Benzo(k)fluoranthene	14.64	252	19886m	3.47	ng	
89) Benzo(a)pyrene	14.93	252	27707m	4.96	ng	
91) Benzo(g,h,i)perylene	16.53	276	17222	3.98	ng	93

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

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