

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011516\
 Data File : BF084511.D
 Acq On : 16 Jan 2016 1:09
 Operator : SJ/UM
 Sample : H1128-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Composite

Quant Time: Jan 18 10:01:26 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 16 03:20:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	252785	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	1042299	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	455962	20.00	ng	-0.01
63) Phenanthrene-d10	11.34	188	807211	20.00	ng	0.00
75) Chrysene-d12	13.99	240	618980	20.00	ng	0.00
86) Perylene-d12	15.40	264	531042	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1887210	120.75	ng	0.01
7) Phenol-d6	6.46	99	2201021	112.14	ng	-0.01
23) Nitrobenzene-d5	7.39	82	1607694	85.85	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	569689	123.76	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	2070538	79.47	ng	-0.01
78) Terphenyl-d14	12.93	244	1634568	58.95	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.58	163	538438	16.52	ng	# 90
70) Phenanthrene	11.37	178	227525	5.39	ng	99
74) Fluoranthene	12.56	202	330667	7.50	ng	97
77) Pyrene	12.79	202	311531	6.41	ng	99
80) Benzo(a)anthracene	13.97	228	118583m	3.26	ng	
82) Chrysene	14.01	228	112766m	3.23	ng	
87) Benzo(b)fluoranthene	15.00	252	128538m	3.70	ng	
89) Benzo(a)pyrene	15.35	252	87001	2.95	ng	96

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

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