

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031517\
 Data File : BF093679.D
 Acq On : 15 Mar 2017 20:26
 Operator : SJ/MA
 Sample : I2264-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-DD8-SSW

Quant Time: Mar 16 03:06:37 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 15:05:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	171111	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	681460	20.00	ng	-0.01
38) Acenaphthene-d10	9.75	164	313121	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	535808	20.00	ng	0.00
75) Chrysene-d12	13.89	240	336674	20.00	ng	0.00
86) Perylene-d12	15.29	264	315020	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1274053	123.92	ng	-0.01
7) Phenol-d6	6.37	99	1621594	125.08	ng	-0.01
23) Nitrobenzene-d5	7.28	82	1024673	83.43	ng	-0.02
41) 2,4,6-Tribromophenol	10.55	330	249660	122.38	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1458813	86.66	ng	-0.01
78) Terphenyl-d14	12.84	244	1000192	77.24	ng	0.01
Target Compounds						
10) Phenol	6.38	94	134882	8.49	ng	Qvalue 89
49) Dimethylphthalate	9.49	163	332391	16.01	ng	99
57) Fluorene	10.31	166	38380	2.13	ng	95
70) Phenanthrene	11.27	178	611987	20.01	ng	99
71) Anthracene	11.33	178	159836	5.17	ng	98
74) Fluoranthene	12.46	202	750062	25.39	ng	96
77) Pyrene	12.69	202	667931	27.28	ng	99
80) Benzo(a)anthracene	13.88	228	289338m	14.55	ng	
82) Chrysene	13.92	228	224232	12.19	ng	99
85) Indeno(1,2,3-cd)pyrene	16.65	276	158918	10.32	ng	96
87) Benzo(b)fluoranthene	14.91	252	295976m	15.43	ng	
88) Benzo(k)fluoranthene	14.92	252	109992m	5.94	ng	
89) Benzo(a)pyrene	15.24	252	244328	13.93	ng	# 94
90) Dibenzo(a,h)anthracene	16.65	278	37624m	2.59	ng	
91) Benzo(a,h,i)perylene	17.05	276	161068	10.82	ng	# 88

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

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