

Data Path : Z:\HPCHEM1\BNA F\DATA\BF063017\
 Data File : BF096362.D
 Acq On : 1 Jul 2017 3:20
 Operator : SJ/MA
 Sample : I3930-11
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-6-8.5

Manual Integrations
 APPROVED

mohammad
 7/3/2017 11:37:53 AM

Quant Time: Jul 01 04:26:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 28 15:41:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	160035	20.00	ng	-0.02
21) Naphthalene-d8	8.04	136	629777	20.00	ng	-0.02
38) Acenaphthene-d10	9.79	164	232564	20.00	ng	-0.02
63) Phenanthrene-d10	11.27	188	355098	20.00	ng	-0.01
75) Chrysene-d12	13.90	240	291786	20.00	ng	-0.01
86) Perylene-d12	15.33	264	199604	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1202020	108.21	ng	0.00
7) Phenol-d6	6.40	99	1446096	107.51	ng	0.00
23) Nitrobenzene-d5	7.32	82	822375	90.40	ng	-0.04
41) 2,4,6-Tribromophenol	10.58	330	251349	137.33	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	1103228	86.88	ng	-0.03
78) Terphenyl-d14	12.86	244	759149	64.92	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	33505	2.27	ng	Qvalue 89
31) Naphthalene	8.06	128	68159	2.25	ng	99
37) 2-Methylnaphthalene	8.75	142	48688	2.47	ng	95
49) Dimethylphthalate	9.51	163	233512	14.03	ng	98
74) Fluoranthene	12.48	202	123974	6.28	ng	96
77) Pyrene	12.71	202	125423	5.23	ng	97
80) Benzo(a)anthracene	13.89	228	52128	2.93	ng	93
82) Chrysene	13.93	228	52368	2.85	ng	97
87) Benzo(b)fluoranthene	14.92	252	58531m	4.46	ng	
89) Benzo(a)pyrene	15.26	252	30101	2.60	ng	# 73
91) Benzo(g,h,i)perylene	17.12	276	22560	2.39	ng	# 87

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

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