

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061317\
 Data File : BF095906.D
 Acq On : 13 Jun 2017 21:50
 Operator : SJ/MA
 Sample : I3629-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CLINTON-AVE-SP-COMP-1

Manual Integrations
 APPROVED

mohammad
 6/14/2017 4:24:47 PM

Quant Time: Jun 14 05:08:55 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.35	152	76808	20.00	ng	-0.05
21) Naphthalene-d8	9.38	136	297363	20.00	ng	-0.05
38) Acenaphthene-d10	12.20	164	119708	20.00	ng	-0.06
63) Phenanthrene-d10	14.60	188	180955	20.00	ng	-0.05
75) Chrysene-d12	18.32	240	154986	20.00	ng	-0.04
86) Perylene-d12	19.99	264	115434	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	481464	99.88	ng	-0.04
7) Phenol-d6	6.92	99	576753	99.95	ng	-0.04
23) Nitrobenzene-d5	8.27	82	326813	68.26	ng	-0.05
41) 2,4,6-Tribromophenol	13.50	330	91096	86.01	ng	-0.05
44) 2-Fluorobiphenyl	11.16	172	560645	65.17	ng	-0.05
78) Terphenyl-d14	16.97	244	333589	53.42	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	11.86	163	63316	6.97	ng	100
70) Phenanthrene	14.64	178	37636	3.60	ng	95
74) Fluoranthene	16.42	202	53285	5.16	ng	98
77) Pyrene	16.72	202	73474	6.35	ng	97
80) Benzo(a)anthracene	18.31	228	27834m	3.09	ng	
82) Chrysene	18.35	228	34830	3.87	ng	93
87) Benzo(b)fluoranthene	19.59	252	24832m	3.44	ng	
89) Benzo(a)pyrene	19.94	252	21906	3.30	ng	# 95
91) Benzo(g,h,i)perylene	21.52	276	12702	2.21	ng	97

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

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