

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061617\
 Data File : BF095992.D
 Acq On : 17 Jun 2017 11:59
 Operator : SJ/MA
 Sample : I3686-04
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-R5-BASE

Manual Integrations
 APPROVED

mohammad
 6/19/2017 3:48:15 PM

Quant Time: Jun 19 01:54:01 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 17:12:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	75330	20.00	ng	0.00
21) Naphthalene-d8	9.35	136	302237	20.00	ng	0.00
38) Acenaphthene-d10	12.17	164	127759	20.00	ng	-0.01
63) Phenanthrene-d10	14.56	188	190936	20.00	ng	0.00
75) Chrysene-d12	18.29	240	148054	20.00	ng	-0.01
86) Perylene-d12	19.96	264	139685	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	436191	92.27	ng	0.00
7) Phenol-d6	6.89	99	550023	97.18	ng	-0.02
23) Nitrobenzene-d5	8.24	82	308528	63.40	ng	-0.01
41) 2,4,6-Tribromophenol	13.47	330	90299	79.88	ng	-0.01
44) 2-Fluorobiphenyl	11.13	172	483266	52.64	ng	-0.01
78) Terphenyl-d14	16.94	244	282597	47.37	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	11.83	163	75120	7.74	ng	100
70) Phenanthrene	14.60	178	119182	10.82	ng	97
71) Anthracene	14.70	178	27763m	2.41	ng	
74) Fluoranthene	16.39	202	147759	13.55	ng	98
77) Pyrene	16.69	202	136643	12.37	ng	96
80) Benzo(a)anthracene	18.27	228	57283	6.65	ng	98
82) Chrysene	18.32	228	54557	6.34	ng	98
85) Indeno(1,2,3-cd)pyrene	21.14	276	22661	3.08	ng	96
87) Benzo(b)fluoranthene	19.56	252	64045	7.32	ng	# 96
88) Benzo(k)fluoranthene	19.59	252	21588m	2.58	ng	
89) Benzo(a)pyrene	19.90	252	50805	6.32	ng	98
91) Benzo(a,h,i)perylene	21.47	276	19800	2.85	ng	97

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

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