

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
 Data File : BF096109.D
 Acq On : 22 Jun 2017 6:24
 Operator : SJ/MA
 Sample : I3747-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-R5-WSW

Manual Integrations
 APPROVED

mohammad
 6/22/2017 6:03:00 PM

Quant Time: Jun 22 07:39:23 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 16:05:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	80912	20.00	ng	0.00
21) Naphthalene-d8	9.32	136	311414	20.00	ng	0.00
38) Acenaphthene-d10	12.15	164	124601	20.00	ng	0.00
63) Phenanthrene-d10	14.54	188	181215	20.00	ng	0.00
75) Chrysene-d12	18.27	240	171796	20.00	ng	0.00
86) Perylene-d12	19.94	264	119012	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	796498	156.86	ng	0.00
7) Phenol-d6	6.87	99	997182	164.04	ng	0.00
23) Nitrobenzene-d5	8.22	82	517738	103.25	ng	0.00
41) 2,4,6-Tribromophenol	13.44	330	129847	117.78	ng	0.00
44) 2-Fluorobiphenyl	11.10	172	846443	94.53	ng	0.00
78) Terphenyl-d14	16.92	244	637765	92.13	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	11.80	163	145777	15.41	ng	99
51) Acenaphthene	12.20	154	32658	4.12	ng	98
57) Fluorene	13.04	166	28458	2.93	ng	# 98
70) Phenanthrene	14.57	178	292057	27.93	ng	98
71) Anthracene	14.67	178	98142	8.99	ng	95
74) Fluoranthene	16.37	202	590437	57.05	ng	99
80) Benzo(a)anthracene	18.26	228	226918	22.72	ng	98
82) Chrysene	18.30	228	171225	17.15	ng	98
85) Indeno(1,2,3-cd)pyrene	21.11	276	67529	7.91	ng	# 90
87) Benzo(b)fluoranthene	19.54	252	169758m	22.78	ng	
88) Benzo(k)fluoranthene	19.57	252	57574m	8.08	ng	
89) Benzo(a)pyrene	19.89	252	132472	19.34	ng	96
90) Dibenzo(a,h)anthracene	21.10	278	17349	2.99	ng	# 76
91) Benzo(g,h,i)perylene	21.44	276	56210	9.49	ng	96

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

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