

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030415\
 Data File : BF077607.D
 Acq On : 5 Mar 2015 5:56
 Operator : TP/IZ
 Sample : G1424-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB17

Manual Integrations
 APPROVED

mohammad
 3/5/2015 2:11:28 PM

Quant Time: Mar 05 11:43:13 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 13:27:09 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.22	152	65614	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	265801	20.00	ng	0.00
38) Acenaphthene-d10	10.97	164	128304	20.00	ng	0.00
63) Phenanthrene-d10	12.81	188	247414	20.00	ng	0.00
75) Chrysene-d12	16.10	240	286178	20.00	ng	-0.01
86) Perylene-d12	17.82	264	257721	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	207990	52.92	ng	0.00
7) Phenol-d6	6.79	99	381071	75.41	ng	0.00
23) Nitrobenzene-d5	7.91	82	241400	56.48	ng	0.00
41) 2,4,6-Tribromophenol	11.95	330	13165	10.66	ng	0.00
44) 2-Fluorobiphenyl	10.15	172	495544	60.22	ng	0.00
78) Terphenyl-d14	14.81	244	555962	45.42	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.65	163	33100	3.67	ng	100
70) Phenanthrene	12.84	178	59268	4.13	ng	99
74) Fluoranthene	14.32	202	212022	13.54	ng	98
77) Pyrene	14.60	202	204646	11.69	ng	99
80) Benzo(a)anthracene	16.09	228	101322	6.04	ng	95
82) Chrysene	16.12	228	98995m	6.29	ng	
85) Indeno(1,2,3-cd)pyrene	19.25	276	57736	3.22	ng	99
87) Benzo(b)fluoranthene	17.38	252	116302m	7.76	ng	
88) Benzo(k)fluoranthene	17.40	252	46322m	3.15	ng	
89) Benzo(a)pyrene	17.75	252	85369	5.98	ng	# 92
91) Benzo(g,h,i)perylene	19.67	276	58696	4.27	ng	97

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

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