

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030215\
 Data File : BF077522.D
 Acq On : 2 Mar 2015 15:08
 Operator : TP/IZ
 Sample : G1399-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB10

Manual Integrations
 APPROVED

apatel
 3/3/2015 6:26:45 PM

Quant Time: Mar 03 05:44:02 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 23:46:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.26	152	94856	20.00	ng	0.00
21) Naphthalene-d8	8.83	136	383033	20.00	ng	0.00
38) Acenaphthene-d10	11.01	164	201924	20.00	ng	0.00
63) Phenanthrene-d10	12.85	188	392443	20.00	ng	0.00
75) Chrysene-d12	16.15	240	406901	20.00	ng	0.02
86) Perylene-d12	17.96	264	411446	20.00	ng	0.13
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	547838	96.42	ng	0.00
7) Phenol-d6	6.82	99	727855	99.63	ng	0.00
23) Nitrobenzene-d5	7.95	82	429851	69.79	ng	0.00
41) 2,4,6-Tribromophenol	11.99	330	180938	93.06	ng	0.00
44) 2-Fluorobiphenyl	10.18	172	861557	66.53	ng	0.00
78) Terphenyl-d14	14.85	244	976238	56.09	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.68	163	32947	2.32	ng	99
70) Phenanthrene	12.88	178	192443	8.46	ng	99
74) Fluoranthene	14.36	202	337745	13.59	ng	96
77) Pyrene	14.64	202	373023	14.99	ng	99
80) Benzo(a)anthracene	16.14	228	195648	8.20	ng	97
82) Chrysene	16.18	228	194338	8.68	ng	97
85) Indeno(1,2,3-cd)pyrene	19.41	276	136813m	5.36	ng	
87) Benzo(b)fluoranthene	17.49	252	246276m	10.29	ng	
88) Benzo(k)fluoranthene	17.52	252	97764m	4.17	ng	
89) Benzo(a)pyrene	17.89	252	206046m	9.04	ng	
91) Benzo(g,h,i)perylene	19.85	276	143456	6.54	ng	96

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

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