

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042115\
 Data File : BF078693.D
 Acq On : 21 Apr 2015 20:05
 Operator : TP/IZ
 Sample : G1938-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-66S

Manual Integrations
 APPROVED

apatel
 4/22/2015 5:22:25 PM

Quant Time: Apr 22 03:33:05 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 00:52:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.82	152	35693	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	148865	20.00	ng	-0.01
38) Acenaphthene-d10	11.17	164	76190	20.00	ng	-0.01
63) Phenanthrene-d10	13.90	188	152888	20.00	ng	-0.01
75) Chrysene-d12	17.77	240	161839	20.00	ng	-0.01
86) Perylene-d12	19.52	264	159287	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	3.84	112	178340	82.38	ng	0.01
7) Phenol-d6	5.32	99	232131	85.56	ng	0.00
23) Nitrobenzene-d5	6.75	82	135919	55.34	ng	-0.01
41) 2,4,6-Tribromophenol	12.64	330	57989	91.77	ng	-0.01
44) 2-Fluorobiphenyl	9.99	172	285043	53.48	ng	-0.01
78) Terphenyl-d14	16.43	244	330242	46.11	ng	0.00
Target Compounds						
49) Dimethylphthalate	10.80	163	13871	2.42	ng	Qvalue # 98
70) Phenanthrene	13.94	178	52314	5.92	ng	97
74) Fluoranthene	15.83	202	83399	8.94	ng	97
77) Pyrene	16.14	202	79383	7.81	ng	96
80) Benzo(a)anthracene	17.76	228	48466	5.03	ng	99
82) Chrysene	17.81	228	35981	3.98	ng	98
85) Indeno(1,2,3-cd)pyrene	20.72	276	28701m	2.80	ng	
87) Benzo(b)fluoranthene	19.07	252	54397m	5.53	ng	
89) Benzo(a)pyrene	19.45	252	39236	4.57	ng	# 92
91) Benzo(g,h,i)perylene	21.04	276	26250	3.04	ng	97

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

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