

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033015\
 Data File : BF078103.D
 Acq On : 31 Mar 2015 4:38
 Operator : TP/IZ
 Sample : G1667-21RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 REC-7-1-4(1.5-2.0)RE

Manual Integrations
 APPROVED

apatel
 3/31/2015 2:17:35 PM

Quant Time: Mar 31 06:45:04 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 28 05:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.06	152	42341	20.00	ng	-0.01
21) Naphthalene-d8	8.64	136	174774	20.00	ng	-0.01
38) Acenaphthene-d10	10.81	164	84144	20.00	ng	-0.01
63) Phenanthrene-d10	12.65	188	169375	20.00	ng	-0.01
75) Chrysene-d12	15.93	240	191742	20.00	ng	0.00
86) Perylene-d12	17.69	264	190453	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	48303	19.91	ng	-0.02
7) Phenol-d6	6.65	99	55305	18.45	ng	-0.02
23) Nitrobenzene-d5	7.76	82	38702	14.52	ng	-0.02
41) 2,4,6-Tribromophenol	11.79	330	27432	34.07	ng	-0.01
44) 2-Fluorobiphenyl	9.98	172	124957	21.82	ng	-0.01
78) Terphenyl-d14	14.64	244	278947	35.54	ng	-0.01
Target Compounds						Qvalue
70) Phenanthrene	12.67	178	22495	2.31	ng	99
74) Fluoranthene	14.15	202	58804	5.48	ng	96
77) Pyrene	14.42	202	56073	4.65	ng	# 95
80) Benzo(a)anthracene	15.92	228	35891	3.16	ng	99
82) Chrysene	15.96	228	30822	3.02	ng	97
87) Benzo(b)fluoranthene	17.23	252	39939m	3.21	ng	
89) Benzo(a)pyrene	17.62	252	29783	2.87	ng	# 97

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

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