

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
 Data File : BF079297.D
 Acq On : 16 May 2015 6:45
 Operator : TP/IZ
 Sample : G2215-14
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-37D

Manual Integrations
 APPROVED

apatel
 5/18/2015 3:42:05 PM

Quant Time: May 18 01:16:18 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 18 00:56:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.18	152	75351	20.00	ng	0.00
21) Naphthalene-d8	8.75	136	303014	20.00	ng	0.00
38) Acenaphthene-d10	10.92	164	140512	20.00	ng	0.00
63) Phenanthrene-d10	12.77	188	260918	20.00	ng	0.00
75) Chrysene-d12	16.06	240	268895	20.00	ng	0.00
86) Perylene-d12	17.78	264	275194	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	417506m	95.38	ng	0.01
7) Phenol-d6	6.75	99	532434	92.02	ng	0.01
23) Nitrobenzene-d5	7.87	82	318886	60.48	ng	0.00
41) 2,4,6-Tribromophenol	11.91	330	144083	94.79	ng	0.00
44) 2-Fluorobiphenyl	10.10	172	566853	56.20	ng	0.00
78) Terphenyl-d14	14.77	244	493342	43.93	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.60	163	33442	3.24	ng	98
70) Phenanthrene	12.80	178	62702	4.30	ng	99
74) Fluoranthene	14.27	202	93956	6.18	ng	92
77) Pyrene	14.55	202	87327	5.64	ng	98
80) Benzo(a)anthracene	16.05	228	49434	3.36	ng	96
82) Chrysene	16.08	228	43545	3.07	ng	97
87) Benzo(b)fluoranthene	17.34	252	64440m	4.28	ng	
89) Benzo(a)pyrene	17.71	252	44369	3.20	ng	# 90
91) Benzo(g,h,i)perylene	19.59	276	35343	2.39	ng	# 94

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

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