

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052215\
 Data File : BF079455.D
 Acq On : 23 May 2015 4:33
 Operator : TP/IZ
 Sample : G2359-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-27D

Manual Integrations
 APPROVED

apatel
 5/27/2015 8:17:13 AM

Quant Time: May 23 06:42:54 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 23 06:12:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.05	152	64319	20.00	ng	0.00
21) Naphthalene-d8	8.63	136	266672	20.00	ng	0.00
38) Acenaphthene-d10	10.80	164	125590	20.00	ng	0.00
63) Phenanthrene-d10	12.63	188	244364	20.00	ng	0.00
75) Chrysene-d12	15.91	240	237732	20.00	ng	0.00
86) Perylene-d12	17.60	264	248780	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	353432m	94.59	ng	0.00
7) Phenol-d6	6.63	99	472174m	95.60	ng	0.00
23) Nitrobenzene-d5	7.75	82	258786	55.77	ng	0.00
41) 2,4,6-Tribromophenol	11.77	330	127928	94.16	ng	0.00
44) 2-Fluorobiphenyl	9.98	172	470835	52.23	ng	0.00
78) Terphenyl-d14	14.63	244	482800	48.62	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.65	128	77220	6.08	ng	99
37) 2-Methylnaphthalene	9.52	142	26653	3.03	ng	99
49) Dimethylphthalate	10.49	163	33139	3.60	ng	# 99
51) Acenaphthene	10.83	154	54452	7.14	ng	100
54) Dibenzofuran	11.05	168	68183	6.19	ng	99
57) Fluorene	11.47	166	62471	6.91	ng	98
70) Phenanthrene	12.66	178	783416	57.31	ng	99
71) Anthracene	12.72	178	203015	15.14	ng	97
72) Carbazole	12.94	167	99031	7.75	ng	98
74) Fluoranthene	14.14	202	725428	50.92	ng	96
77) Pyrene	14.41	202	622265	45.42	ng	98
80) Benzo(a)anthracene	15.90	228	328889	25.26	ng	98
82) Chrysene	15.94	228	254640	20.34	ng	97
85) Indeno(1,2,3-cd)pyrene	18.94	276	153798	9.64	ng	98
87) Benzo(b)fluoranthene	17.18	252	312963m	22.98	ng	
88) Benzo(k)fluoranthene	17.20	252	130194m	9.88	ng	
89) Benzo(a)pyrene	17.54	252	259471	20.69	ng	98
90) Dibenzo(a,h)anthracene	18.95	278	41663	3.13	ng	# 90
91) Benzo(a,h,i)perylene	19.33	276	153266	11.48	ng	98

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

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