

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
 Data File : BF079294.D
 Acq On : 16 May 2015 5:19
 Operator : TP/IZ
 Sample : G2259-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WASTE-1

Manual Integrations
 APPROVED

apatel
 5/18/2015 3:41:47 PM

Quant Time: May 16 06:27:43 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 16 04:02:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.18	152	66841	20.00	ng	0.00
21) Naphthalene-d8	8.75	136	275026	20.00	ng	0.00
38) Acenaphthene-d10	10.93	164	129021	20.00	ng	0.00
63) Phenanthrene-d10	12.77	188	229254	20.00	ng	0.00
75) Chrysene-d12	16.06	240	224694	20.00	ng	0.00
86) Perylene-d12	17.76	264	246065	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	513321m	132.20	ng	0.00
7) Phenol-d6	6.75	99	682837	133.04	ng	0.01
23) Nitrobenzene-d5	7.87	82	407042	85.06	ng	0.00
41) 2,4,6-Tribromophenol	11.91	330	179963	128.93	ng	0.00
44) 2-Fluorobiphenyl	10.10	172	733937	79.25	ng	0.00
78) Terphenyl-d14	14.77	244	654746	69.76	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.78	128	32775	2.50	ng	98
48) Acenaphthylene	10.75	152	48933	3.73	ng	97
49) Dimethylphthalate	10.61	163	89423	9.44	ng	99
54) Dibenzofuran	11.18	168	30714	2.71	ng	97
57) Fluorene	11.60	166	35915	3.87	ng	98
70) Phenanthrene	12.80	178	699519	54.54	ng	99
71) Anthracene	12.86	178	132323	10.52	ng	98
72) Carbazole	13.06	167	43949	3.66	ng	99
74) Fluoranthene	14.27	202	894049	66.90	ng	98
77) Pyrene	14.55	202	831153	64.19	ng	97
80) Benzo(a)anthracene	16.03	228	398607m	32.40	ng	
82) Chrysene	16.08	228	357942	30.25	ng	98
83) Bis(2-ethylhexyl)phthalate	16.10	149	71351	8.32	ng	99
85) Indeno(1,2,3-cd)pyrene	19.15	276	193523	12.83	ng	94
87) Benzo(b)fluoranthene	17.33	252	424327m	31.51	ng	
88) Benzo(k)fluoranthene	17.35	252	166148m	12.74	ng	
89) Benzo(a)pyrene	17.69	252	329713	26.59	ng	97
90) Dibenzo(a,h)anthracene	19.17	278	45207m	3.43	ng	
91) Benzo(a,h,i)perylene	19.57	276	177772	13.46	ng	97

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

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