

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090815\
 Data File : BF081527.D
 Acq On : 8 Sep 2015 17:39
 Operator : UM/IZ
 Sample : G3472-01DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB001DL

Manual Integrations
 APPROVED

MMDadoda
 9/9/2015 2:58:13 PM

Quant Time: Sep 09 01:19:23 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 23:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.34	152	46515	20.00	ng	0.00
21) Naphthalene-d8	7.63	136	187361	20.00	ng	-0.01
38) Acenaphthene-d10	9.37	164	85522	20.00	ng	-0.01
63) Phenanthrene-d10	10.83	188	161001	20.00	ng	-0.01
75) Chrysene-d12	13.45	240	126361	20.00	ng	0.00
86) Perylene-d12	14.75	264	89822	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.88	112	56921	18.99	ng	0.00
7) Phenol-d6	6.01	99	80738	20.35	ng	-0.01
23) Nitrobenzene-d5	6.92	82	46188	11.89	ng	0.00
41) 2,4,6-Tribromophenol	10.16	330	13738	18.04	ng	0.00
44) 2-Fluorobiphenyl	8.72	172	87782	13.15	ng	0.00
78) Terphenyl-d14	12.42	244	73965	13.63	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.12	163	17258	2.60	ng	# 95
57) Fluorene	9.92	166	13851	2.18	ng	97
70) Phenanthrene	10.86	178	184078	20.57	ng	98
71) Anthracene	10.91	178	47285	5.14	ng	99
74) Fluoranthene	12.03	202	308606	31.85	ng	91
77) Pyrene	12.26	202	281004	30.02	ng	99
80) Benzo(a)anthracene	13.44	228	131004m	16.28	ng	
82) Chrysene	13.47	228	107907m	14.96	ng	
85) Indeno(1,2,3-cd)pyrene	15.82	276	51320	9.70	ng	# 83
87) Benzo(b)fluoranthene	14.43	252	96903m	15.11	ng	
88) Benzo(k)fluoranthene	14.43	252	48926m	9.19	ng	
89) Benzo(a)pyrene	14.71	252	78656	14.47	ng	# 83
90) Dibenzo(a,h)anthracene	15.83	278	11857	2.82	ng	# 81
91) Benzo(g,h,i)perylene	16.14	276	47477	11.85	ng	# 72

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

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