

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082115\
 Data File : BF081157.D
 Acq On : 21 Aug 2015 21:19
 Operator : UM/IZ
 Sample : G3363-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-C1

Manual Integrations
 APPROVED

MMDadoda
 8/24/2015 11:24:02 AM

Quant Time: Aug 22 05:34:15 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 21 23:26:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	68081	20.00	ng	-0.01
21) Naphthalene-d8	7.84	136	294363	20.00	ng	0.00
38) Acenaphthene-d10	9.58	164	124112	20.00	ng	-0.01
63) Phenanthrene-d10	11.05	188	265226	20.00	ng	0.00
75) Chrysene-d12	13.67	240	181331	20.00	ng	0.00
86) Perylene-d12	15.01	264	153567	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	499645	110.39	ng	0.02
7) Phenol-d6	6.21	99	625807	105.97	ng	0.00
23) Nitrobenzene-d5	7.12	82	391201	60.71	ng	-0.01
41) 2,4,6-Tribromophenol	10.37	330	106411	98.91	ng	0.00
44) 2-Fluorobiphenyl	8.92	172	517596	54.64	ng	-0.01
78) Terphenyl-d14	12.64	244	477396	56.44	ng	0.01
Target Compounds						Qvalue
31) Naphthalene	7.86	128	35357	2.15	ng	99
49) Dimethylphthalate	9.32	163	121546	11.89	ng	100
51) Acenaphthene	9.61	154	37479	3.98	ng	99
54) Dibenzofuran	9.78	168	41517	3.29	ng	99
57) Fluorene	10.13	166	39323	4.06	ng	98
70) Phenanthrene	11.08	178	480130	30.55	ng	100
71) Anthracene	11.12	178	147331	9.63	ng	98
72) Carbazole	11.28	167	72409	4.92	ng	99
74) Fluoranthene	12.25	202	609628	35.94	ng	92
77) Pyrene	12.48	202	603289	40.93	ng	100
80) Benzo(a)anthracene	13.66	228	250758m	20.62	ng	
82) Chrysene	13.69	228	222880m	20.34	ng	
83) Bis(2-ethylhexyl)phthalate	13.68	149	18640	2.30	ng	97
85) Indeno(1,2,3-cd)pyrene	16.19	276	102620	13.34	ng	95
87) Benzo(b)fluoranthene	14.65	252	240583m	22.15	ng	
88) Benzo(k)fluoranthene	14.67	252	67421m	6.66	ng	
89) Benzo(a)pyrene	14.95	252	169185	17.88	ng	# 94
90) Dibenzo(a,h)anthracene	16.19	278	26189m	3.55	ng	
91) Benzo(g,h,i)perylene	16.54	276	85683	11.45	ng	# 92

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

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