

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
 Data File : BE090009.D
 Acq On : 30 Jun 2015 8:54
 Operator : TP/IZ
 Sample : G2811-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0520624-150625

Manual Integrations
 APPROVED

apatel
 6/30/2015 4:55:42 PM

Quant Time: Jun 30 16:29:31 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	16540	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	81449	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	48642	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	109467	5.00	ng	0.00
25) Chrysene-d12	21.13	240	115299	5.00	ng	0.00
32) Perylene-d12	23.46	264	116708	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	34433	9.06	ng	0.00
5) Phenol-d6	6.81	99	56282	10.58	ng	0.00
7) Nitrobenzene-d5	8.76	82	33842	5.23	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	18331	12.86	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	72090	4.93	ng	0.00
27) Terphenyl-d14	19.59	244	103822	5.40	ng	0.00
Target Compounds						
9) Naphthalene	10.43	128	823	0.04	ng	# 74
11) 2-Methylnaphthalene	12.05	142	262	0.02	ng	# 85
16) Acenaphthene	14.30	154	657	0.05	ng	98
17) Fluorene	15.29	166	650	0.04	ng	# 88
22) Phenanthrene	17.01	178	16484	0.59	ng	99
23) Anthracene	17.10	178	3355m	0.13	ng	
24) Fluoranthene	19.02	202	54644	1.74	ng	99
26) Pyrene	19.38	202	49955	1.82	ng	# 95
28) Benzo(a)anthracene	21.11	228	37191	1.29	ng	97
29) Chrysene	21.17	228	38681	1.29	ng	96
31) Indeno(1,2,3-cd)pyrene	25.87	276	26955	0.94	ng	95
33) Benzo(b)fluoranthene	22.75	252	54927	2.13	ng	# 95
34) Benzo(k)fluoranthene	22.79	252	19233m	0.67	ng	
35) Benzo(a)pyrene	23.35	252	37642	1.58	ng	96
36) Dibenzo(a,h)anthracene	25.89	278	6675	0.27	ng	# 80
37) Benzo(g,h,i)perylene	26.62	276	27544	1.11	ng	96

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

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