

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081215\
 Data File : BE090490.D
 Acq On : 13 Aug 2015 5:01
 Operator : UM/IZ
 Sample : G3248-02
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 CSB-01

Manual Integrations
 APPROVED

MMDadoda
 8/13/2015 3:35:22 PM

Quant Time: Aug 13 13:34:39 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:28:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	20105	5.00	ng	0.00
6) Naphthalene-d8	10.34	136	94496	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	53957	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	125968	5.00	ng	0.00
25) Chrysene-d12	21.09	240	149307	5.00	ng	0.00
32) Perylene-d12	23.39	264	141822	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0d	0.00	ng	
5) Phenol-d6	6.77	99	2052	0.55	ng	0.00
7) Nitrobenzene-d5	8.72	82	25123	3.34	ng	-0.01
13) 2,4,6-Tribromophenol	0.00	330	0d	0.00	ng	
14) 2-Fluorobiphenyl	12.82	172	57927	3.90	ng	0.00
27) Terphenyl-d14	19.55	244	94794	4.31	ng	-0.01
Target Compounds						
9) Naphthalene	10.38	128	923	0.04	ng	# 79
11) 2-Methylnaphthalene	12.01	142	397	0.03	ng	# 89
22) Phenanthrene	16.98	178	6935	0.22	ng	96
24) Fluoranthene	18.98	202	12678	0.41	ng	99
26) Pyrene	19.34	202	7822	0.24	ng	98
28) Benzo(a)anthracene	21.08	228	3793	0.11	ng	# 91
29) Chrysene	21.13	228	6044	0.15	ng	# 92
30) Bis(2-ethylhexyl)phthalate	21.03	149	3532	0.54	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	3194	0.34	ng	99
33) Benzo(b)fluoranthene	22.70	252	6827m	0.40	ng	
34) Benzo(k)fluoranthene	22.74	252	2659m	0.07	ng	
37) Benzo(g,h,i)perylene	26.51	276	3273	0.12	ng	# 88

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

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