

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041116\
 Data File : BE091637.D
 Acq On : 11 Apr 2016 20:32
 Operator : UM/SJ
 Sample : H2248-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TILCONMHDGA005

Manual Integrations
 APPROVED

Sohil
 4/12/2016 5:35:10 PM

Quant Time: Apr 12 15:45:08 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 16:48:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	41870	5.00	ng	-0.02
7) Naphthalene-d8	10.59	136	204841	5.00	ng	0.00
13) Acenaphthene-d10	14.42	164	126072	5.00	ng	0.00
19) Phenanthrene-d10	17.16	188	318189	5.00	ng	0.00
26) Chrysene-d12	21.31	240	418642	5.00	ng	0.00
35) Perylene-d12	23.76	264	378929	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	36957	3.70	ng	0.00
5) Phenol-d6	6.97	99	64464	4.84	ng	0.00
8) Nitrobenzene-d5	8.96	82	29026	2.44	ng	0.00
14) 2,4,6-Tribromophenol	15.90	330	16860	3.79	ng	-0.01
15) 2-Fluorobiphenyl	13.05	172	99481	2.84	ng	0.00
29) Terphenyl-d14	19.76	244	167811	3.21	ng	0.00
Target Compounds						
22) Pentachlorophenol	16.81	266	101	0.24	ng	# 80
23) Phenanthrene	17.19	178	1908m	0.03	ng	
25) Fluoranthene	19.21	202	2376	0.03	ng	# 87
28) Pyrene	19.57	202	2738	0.03	ng	# 91
30) Benzo(a)anthracene	21.29	228	3776m	0.04	ng	
32) Chrysene	21.34	228	2708m	0.03	ng	
34) Indeno(1,2,3-cd)pyrene	26.32	276	2661	0.03	ng	94
40) Benzo(g,h,i)perylene	27.11	276	2834	0.03	ng	# 64

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

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