

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112215\
 Data File : BE091288.D
 Acq On : 21 Nov 2015 22:19
 Operator : UM/NP
 Sample : G4345-24
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4J69

Manual Integrations
 APPROVED

Mmdadoda
 11/23/2015 2:14:36 PM

Quant Time: Nov 23 10:35:52 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 22 01:59:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	2838	0.40	ng/ul	-0.02
4) Naphthalene-d8	9.54	136	13352	0.40	ng/ul	-0.02
8) Acenaphthene-d10	13.36	164	7657	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.08	188	17613m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.23	240	19417	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	15168m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.92	96	22867	11.20	ng/uL	-0.03
6) 2-Methylnaphthalene-d10	11.05	152	5462	0.30	ng/ul	0.00
16) Fluoranthene-d10	18.06	212	14163	0.29	ng/ul	0.00
Target Compounds						
10) Acenaphthene	13.42	153	760	0.03	ng/ul	Ovalue 92
11) Fluorene	14.40	166	1246	0.04	ng/ul#	93
13) Pentachlorophenol	15.86	266	199	0.12	ng/ul	94
14) Phenanthrene	16.11	178	65310	0.32	ng/ul	100
17) Fluoranthene	18.09	202	63269	0.27	ng/ul	90
20) Benzo(a)anthracene	20.20	228	5129m	0.10	ng/ul	
21) Chrysene	20.26	228	5599	0.10	ng/ul	96
23) Benzo(b)fluoranthene	21.55	252	3747	0.08	ng/ul	91
24) Benzo(k)fluoranthene	21.58	252	3460m	0.07	ng/ul	
26) Indeno(1,2,3-cd)pyrene	23.71	276	8727	0.05	ng/ul	94
28) Benzo(g,h,i)perylene	24.26	276	10026m	0.05	ng/ul	

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

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