

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112215\
 Data File : BE091289.D
 Acq On : 21 Nov 2015 22:55
 Operator : UM/NP
 Sample : G4345-26MS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4J69MS

Manual Integrations
 APPROVED

Mmdadoda
 11/23/2015 2:14:37 PM

Quant Time: Nov 23 10:36:50 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	6.89	152	2647	0.40	ng/ul	0.00
4) Naphthalene-d8	9.56	136	12735	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.36	164	6950	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.09	188	16682	0.40	ng/ul	0.00
18) Chrysene-d12	20.23	240	19470	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	17233m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	11.06	152	5679	0.33	ng/ul	0.00
16) Fluoranthene-d10	18.06	212	17644	0.38	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	9.60	128	1543	0.05	ng/ul#	84
10) Acenaphthene	13.42	153	1449	0.06	ng/ul	93
11) Fluorene	14.41	166	2247	0.07	ng/ul#	97
13) Pentachlorophenol	15.86	266	899	0.57	ng/ul	96
14) Phenanthrene	16.12	178	120598	0.63	ng/ul	96
15) Anthracene	16.20	178	23129	0.13	ng/ul	96
17) Fluoranthene	18.10	202	119369	0.53	ng/ul	88
19) Pyrene	18.47	202	92422	0.42	ng/ul#	79
20) Benzo(a)anthracene	20.20	228	9310m	0.19	ng/ul	
21) Chrysene	20.26	228	10068	0.17	ng/ul	98
23) Benzo(b)fluoranthene	21.55	252	6994	0.13	ng/ul	97
24) Benzo(k)fluoranthene	21.58	252	6447m	0.12	ng/ul	
26) Indeno(1,2,3-cd)pyrene	23.70	276	17752m	0.09	ng/ul	
28) Benzo(g,h,i)perylene	24.26	276	19958m	0.09	ng/ul	

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

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