

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112215\
 Data File : BE091279.D
 Acq On : 21 Nov 2015 16:06
 Operator : UM/NP
 Sample : G4346-08MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4J72MSD

Quant Time: Nov 23 10:26:24 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 : Mon Nov 23 00:41:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	3188	0.40	ng/ul	-0.02
4) Naphthalene-d8	9.56	136	14023	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.36	164	7823	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.09	188	18618	0.40	ng/ul	0.00
18) Chrysene-d12	20.23	240	22367	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	19932	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	11.05	152	6776	0.35	ng/ul	0.00
16) Fluoranthene-d10	18.07	212	18949	0.36	ng/ul	0.00
Target Compounds						
					Ovalue	
10) Acenaphthene	13.42	153	155	0.01	ng/ul#	75
13) Pentachlorophenol	15.85	266	1509	0.85	ng/ul	97
14) Phenanthrene	16.13	178	17716	0.08	ng/ul#	95
15) Anthracene	16.20	178	3223	0.02	ng/ul#	48
17) Fluoranthene	18.09	202	20303	0.08	ng/ul	87
19) Pyrene	18.48	202	19595	0.08	ng/ul#	78
20) Benzo(a)anthracene	20.20	228	2718	0.05	ng/ul#	93
21) Chrysene	20.26	228	3540	0.05	ng/ul#	91
26) Indeno(1,2,3-cd)pyrene	23.71	276	5082	0.02	ng/ul	98
28) Benzo(g,h,i)perylene	24.26	276	5628	0.02	ng/ul#	42

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

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