

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112115\
 Data File : BE091260.D
 Acq On : 21 Nov 2015 1:26
 Operator : UM/NP
 Sample : G4310-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 D9KN1

Manual Integrations
 APPROVED

Mmdadoda
 11/23/2015 2:13:20 PM

Quant Time: Nov 23 09:46:37 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 : Sat Nov 21 04:53:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	2589	0.40	ng/ul	-0.03
4) Naphthalene-d8	9.55	136	12429	0.40	ng/ul	-0.03
8) Acenaphthene-d10	13.36	164	6929	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.08	188	17345m	0.40	ng/ul	-0.02
18) Chrysene-d12	20.23	240	19570	0.40	ng/ul	0.00
22) Perylene-d12	22.11	264	17642m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.93	96	24653	13.24	ng/uL	-0.04
6) 2-Methylnaphthalene-d10	11.06	152	6255	0.37	ng/ul	-0.02
16) Fluoranthene-d10	18.07	212	19399	0.40	ng/ul	0.00
Target Compounds						
13) Pentachlorophenol	15.85	266	357	0.22	ng/ul	96
14) Phenanthrene	16.11	178	23915	0.12	ng/ul	96
17) Fluoranthene	18.10	202	66119	0.28	ng/ul	88
20) Benzo(a)anthracene	20.20	228	3848m	0.08	ng/ul	
21) Chrysene	20.26	228	12123	0.21	ng/ul	96
23) Benzo(b)fluoranthene	21.55	252	11331	0.21	ng/ul	98
24) Benzo(k)fluoranthene	21.58	252	7464m	0.13	ng/ul	
25) Benzo(a)pyrene	22.01	252	7382	0.15	ng/ul	96
26) Indeno(1,2,3-cd)pyrene	23.71	276	22671m	0.11	ng/ul	
28) Benzo(g,h,i)perylene	24.26	276	27537m	0.12	ng/ul	

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

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