

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111115\
 Data File : BE091144.D
 Acq On : 12 Nov 2015 14:37
 Operator : UM/NP
 Sample : G4322-18
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A4HS9

Manual Integrations
 APPROVED

Mmdadoda
 11/16/2015 12:42:36 PM

Quant Time: Nov 13 17:09:58 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 : Fri Nov 13 13:56:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	3522	0.40	ng/ul	0.00
4) Naphthalene-d8	9.57	136	19470	0.40	ng/ul	0.00
8) Acenaphthene-d10	13.40	164	11788	0.40	ng/ul	0.00
12) Phenanthrene-d10	16.12	188	28893m	0.40	ng/ul	-0.01
18) Chrysene-d12	20.29	240	38540	0.40	ng/ul	0.00
22) Perylene-d12	22.19	264	42234	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	2.92	96	8573	3.38	ng/uL	-0.03
6) 2-Methylnaphthalene-d10	11.07	152	3040	0.11	ng/ul	0.00
16) Fluoranthene-d10	18.12	212	8104	0.10	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	13.14	152	2872	0.05	ng/ul	97
11) Fluorene	14.45	166	1554	0.03	ng/ul#	89
14) Phenanthrene	16.16	178	144497	0.43	ng/ul	98
15) Anthracene	16.25	178	31424m	0.10	ng/ul	
17) Fluoranthene	18.15	202	296925	0.76	ng/ul	90
19) Pyrene	18.53	202	257981	0.59	ng/ul#	79
20) Benzo(a)anthracene	20.26	228	29616	0.30	ng/ul	97
21) Chrysene	20.31	228	37590	0.33	ng/ul	98
23) Benzo(b)fluoranthene	21.62	252	30413	0.23	ng/ul	95
24) Benzo(k)fluoranthene	21.65	252	28332	0.21	ng/ul	95
25) Benzo(a)pyrene	22.09	252	30967	0.25	ng/ul	95
26) Indeno(1,2,3-cd)pyrene	23.83	276	84655m	0.17	ng/ul	
27) Dibenzo(a,h)anthracene	23.79	278	5418	0.04	ng/ul#	70
28) Benzo(g,h,i)perylene	24.39	276	73246	0.14	ng/ul	97

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

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