

Data Path : Z:\HPCHEM1\BNA_E\Data\BE030515\
 Data File : BE089750.D
 Acq On : 5 Mar 2015 19:27
 Operator : TP/JJ
 Sample : SST00.409
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 06 09:46:08 2015

Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM01.2-EPA-SIM-BE030315.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Mar 06 09:36:30 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.33	152	5188	0.40	ng/ul	0.00
4) Naphthalene-d8	12.23	136	25242	0.40	ng/ul	0.00
8) Acenaphthene-d10	16.40	164	9498	0.40	ng/ul	0.00
12) Phenanthrene-d10	19.76	188	16735	0.40	ng/ul	0.00
18) Chrysene-d12	23.65	240	10168	0.40	ng/ul	0.00
22) Perylene-d12	25.33	264	6151	0.40	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	2.63	96	15402	2.26	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.87	152	11053	0.41	ng/ul	0.00
16) Fluoranthene-d10	21.69	212	10353	0.37	ng/ul	0.00

Target Compounds

						Qvalue
3) 1,4-Dioxane	2.69	88	17687	2.24	ng/ul	99
5) Naphthalene	12.28	128	21946	0.38	ng/ul	99
7) 2-Methylnaphthalene	13.94	142	13780	0.40	ng/ul	99
9) Acenaphthylene	16.05	152	69975	0.52	ng/ul	97
10) Acenaphthene	16.47	153	40428	0.40	ng/ul	96
11) Fluorene	17.72	166	11071	0.41	ng/ul	98
13) Pentachlorophenol	19.42	266	230	0.35	ng/ul	95
14) Phenanthrene	19.80	178	63450	0.38	ng/ul	98
15) Anthracene	19.89	178	53667	0.38	ng/ul	100
17) Fluoranthene	21.72	202	58682	0.37	ng/ul	99
19) Pyrene	22.03	202	56781	0.43	ng/ul	100
20) Benzo(a)anthracene	23.64	228	4625	0.33	ng/ul	100
21) Chrysene	23.68	228	12030	0.41	ng/ul	98
23) Benzo(b)fluoranthene	24.90	252	1652	0.26	ng/ul	94
24) Benzo(k)fluoranthene	24.93	252	7456	0.37	ng/ul	97
25) Benzo(a)pyrene	25.27	252	2531	0.31	ng/ul	97
26) Indeno(1,2,3-cd)pyrene	26.72	276	9153	0.31	ng/ul	94
27) Dibenzo(a,h)anthracene	26.76	278	1648	0.31	ng/ul	97
28) Benzo(g,h,i)perylene	27.15	276	17545	0.43	ng/ul	97

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

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