

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071415\
 Data File : BE090099.D
 Acq On : 14 Jul 2015 15:30
 Operator : TP/IZ
 Sample : SST0.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.202

Quant Time: Jul 15 00:45:44 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 15 00:44:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1463	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6652	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3888	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8726	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	8082	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5850	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.96	152	2180	0.21	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	4990	0.21	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	3712	0.19	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	2449	0.19	ng/ul	100
7) Acenaphthylene	13.93	152	16236	0.47	ng/ul	100
8) Acenaphthene	14.28	153	11522	0.47	ng/ul	98
9) Fluorene	15.27	166	3407	0.19	ng/ul	98
11) Pentachlorophenol	16.62	266	141	0.07	ng/ul	96
12) Phenanthrene	16.99	178	21695	0.48	ng/ul	99
13) Anthracene	17.09	178	20570	0.47	ng/ul	98
15) Fluoranthene	19.01	202	22893	0.46	ng/ul	96
17) Pyrene	19.37	202	23272	0.45	ng/ul	99
18) Benzo(a)anthracene	21.10	228	4231	0.14	ng/ul	99
19) Chrysene	21.15	228	5586	0.18	ng/ul	98
21) Benzo(b)fluoranthene	22.73	252	3292	0.13	ng/ul	95
22) Benzo(k)fluoranthene	22.78	252	4730	0.21	ng/ul#	94
23) Benzo(a)pyrene	23.33	252	3295	0.15	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.85	276	14666	0.43	ng/ul	99
25) Dibenzo(a,h)anthracene	25.86	278	2628	0.14	ng/ul	96
26) Benzo(g,h,i)perylene	26.58	276	13528	0.45	ng/ul	96

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

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