

Data Path : Z:\HPCHEM1\BNA E\Data\BE101415\
 Data File : BE091009.D
 Acq On : 14 Oct 2015 18:11
 Operator : UM/IZ
 Sample : SSTD0.475
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.475

Quant Time: Oct 15 05:35:12 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE101415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 05:25:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	6223	0.40	ng/ul	0.00
2) Naphthalene-d8	9.63	136	28033	0.40	ng/ul	0.00
6) Acenaphthene-d10	13.45	164	15673	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.16	188	40261	0.40	ng/ul	0.00
16) Chrysene-d12	20.32	240	54292	0.40	ng/ul	0.00
20) Perylene-d12	22.22	264	55214	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.13	152	12217	0.05	ng/ul	0.00
14) Fluoranthene-d10	18.15	212	36635	0.34	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	9.67	128	24215	0.35	ng/ul	100
5) 2-Methylnaphthalene	11.20	142	13532	0.31	ng/ul	100
7) Acenaphthylene	13.19	152	22217	0.09	ng/ul	100
8) Acenaphthene	13.51	153	16158	0.09	ng/ul	100
9) Fluorene	14.49	166	20186	0.32	ng/ul	100
11) Pentachlorophenol	15.92	266	1677	0.32	ng/ul	99
12) Phenanthrene	16.20	178	138507	0.85	ng/ul	100
13) Anthracene	16.28	178	135407	0.86	ng/ul	100
15) Fluoranthene	18.18	202	173220	0.88	ng/ul	100
17) Pyrene	18.56	202	178143	0.85	ng/ul	100
18) Benzo(a)anthracene	20.28	228	46137	0.32	ng/ul	100
19) Chrysene	20.35	228	49346	0.32	ng/ul	100
21) Benzo(b)fluoranthene	21.65	252	50104	0.30	ng/ul	100
22) Benzo(k)fluoranthene	21.68	252	52602	0.33	ng/ul	100
23) Benzo(a)pyrene	22.12	252	49728	0.32	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	23.87	276	195155	0.82	ng/ul	100
25) Dibenzo(a,h)anthracene	23.83	278	50665	0.31	ng/ul	100
26) Benzo(g,h,i)perylene	24.44	276	209221	0.83	ng/ul	100

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

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