

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020216\
 Data File : BE091379.D
 Acq On : 2 Feb 2016 16:05
 Operator : UM/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.8

Quant Time: Feb 02 17:32:38 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 02 17:28:11 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	44425	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	208219	5.00	ng	0.00
8) Acenaphthene-d10	13.36	164	130873	5.00	ng	0.00
14) Phenanthrene-d10	16.09	188	335823	5.00	ng	0.00
18) Chrysene-d12	20.21	240	372016	5.00	ng	0.00
24) Perylene-d12	22.10	264	343180	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.95	112	5677	1.20	ng	0.00
3) Phenol-d6	6.58	99	8799	1.16	ng	0.00
5) Nitrobenzene-d5	8.18	82	6652	0.55	ng	0.00
9) 2,4,6-Tribromophenol	14.91	330	2358	0.88	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	24547	0.87	ng	0.00
20) Terphenyl-d14	18.62	244	39961	1.13	ng	0.00

Target Compounds				Qvalue		
6) Naphthalene	9.59	128	34094	0.77	ng	99
7) 2-Methylnaphthalene	11.14	142	17434	1.00	ng	91
11) Acenaphthylene	13.10	152	125099	0.73	ng	100
12) Acenaphthene	13.42	154	25709	0.69	ng	# 71
13) Fluorene	14.40	166	33703	0.82	ng	95
15) Phenanthrene	16.12	178	64838	0.78	ng	99
16) Anthracene	16.21	178	52060	0.89	ng	99
17) Fluoranthene	18.10	202	56647	0.94	ng	99
19) Pyrene	18.48	202	57324	1.01	ng	99
21) Benzo(a)anthracene	20.19	228	45696	0.88	ng	98
22) Chrysene	20.25	228	72765	0.81	ng	100
23) Indeno(1,2,3-cd)pyrene	23.70	276	46810	0.79	ng	95
25) Benzo(b)fluoranthene	21.55	252	50062	0.87	ng	# 96
26) Benzo(k)fluoranthene	21.57	252	59878	0.89	ng	# 97
27) Benzo(a)pyrene	22.00	252	43112	0.73	ng	# 96
28) Dibenzo(a,h)anthracene	23.65	278	37849	0.87	ng	96
29) Benzo(g,h,i)perylene	24.25	276	50692	0.91	ng	# 92

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

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