

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020216\
 Data File : BE091384.D
 Acq On : 2 Feb 2016 19:09
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE020216

Quant Time: Feb 03 00:35:50 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 00:27:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	45623	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	223820	5.00	ng	0.00
8) Acenaphthene-d10	13.36	164	146695	5.00	ng	0.00
14) Phenanthrene-d10	16.09	188	371147	5.00	ng	0.00
18) Chrysene-d12	20.22	240	440148	5.00	ng	0.00
24) Perylene-d12	22.10	264	430082	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	4.94	112	8882	1.02	ng	0.00
3) Phenol-d6	6.56	99	14025	1.10	ng	0.00
5) Nitrobenzene-d5	8.16	82	11791	1.03	ng	0.00
9) 2,4,6-Tribromophenol	14.91	330	4460	0.93	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	36241	0.95	ng	0.00
20) Terphenyl-d14	18.62	244	67623	1.04	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	9.59	128	48500	0.96	ng	100
7) 2-Methylnaphthalene	11.13	142	27466	0.98	ng	98
11) Acenaphthylene	13.10	152	202290	0.94	ng	99
12) Acenaphthene	13.42	154	37134	0.92	ng	99
13) Fluorene	14.40	166	50170	0.96	ng	99
15) Phenanthrene	16.12	178	91315	0.94	ng	100
16) Anthracene	16.21	178	80127	0.98	ng	99
17) Fluoranthene	18.09	202	100927	1.08	ng	100
19) Pyrene	18.47	202	98161	1.03	ng	100
21) Benzo(a)anthracene	20.19	228	89446	1.05	ng	98
22) Chrysene	20.25	228	111755	0.97	ng	99
23) Indeno(1,2,3-cd)pyrene	23.69	276	94384	1.07	ng	97
25) Benzo(b)fluoranthene	21.55	252	96234	1.02	ng	99
26) Benzo(k)fluoranthene	21.57	252	102729	0.97	ng	99
27) Benzo(a)pyrene	22.00	252	90290	1.05	ng	99
28) Dibenzo(a,h)anthracene	23.64	278	85646	1.08	ng	99
29) Benzo(g,h,i)perylene	24.24	276	95483	1.01	ng	100

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

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