

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070117\
 Data File : BE093378.D
 Acq On : 30 Jun 2017 15:56
 Operator : SJ/JU
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

mohammad
 7/3/2017 1:14:06 PM

Quant Time: Jun 30 19:59:40 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 30 18:48:18 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	3172	0.40	ng	0.00
7) Naphthalene-d8	10.74	136	15233	0.40	ng	0.00
13) Acenaphthene-d10	14.56	164	9548	0.40	ng	0.00
19) Phenanthrene-d10	17.27	188	38493	0.40	ng	0.00
27) Chrysene-d12	21.44	240	38537	0.40	ng	0.00
34) Perylene-d12	23.95	264	35625	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.53	112	1017	0.11	ng	0.00
5) Phenol-d6	7.10	99	1519	0.11	ng	0.00
8) Nitrobenzene-d5	9.09	82	980	0.08	ng	0.00
11) 2-Methylnaphthalene-d10	12.32	152	2319	0.10	ng	0.00
14) 2,4,6-Tribromophenol	16.03	330	348	0.16	ng	0.00
15) 2-Fluorobiphenyl	13.19	172	3625	0.09	ng	0.01
25) Fluoranthene-d10	19.30	212	34348	0.27	ng	0.00
29) Terphenyl-d14	19.89	244	6950	0.09	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.45	88	725	0.103	ng	# 91
6) bis(2-Chloroethyl)ether	7.36	93	1012	0.085	ng	# 91
9) Naphthalene	10.79	128	15791	0.386	ng	# 74
10) Hexachlorobutadiene	11.08	225	605	0.109	ng	92
12) 2-Methylnaphthalene	12.39	142	2559	0.101	ng	96
16) Acenaphthylene	14.27	152	3865	0.023	ng	# 86
17) Acenaphthene	14.61	154	2800	0.090	ng	97
18) Fluorene	15.59	166	3840	0.103	ng	# 86
20) 4-Bromophenyl-phenylether	16.45	248	729	0.031	ng	# 51
21) Hexachlorobenzene	16.58	284	1207	0.059	ng	# 100
22) Pentachlorophenol	16.94	266	153m	0.023	ng	
23) Phenanthrene	17.30	178	8247	0.047	ng	95
24) Anthracene	17.40	178	9092	0.062	ng	98
28) Pyrene	19.69	202	9983	0.021	ng	# 97
30) Benzo(a)anthracene	21.42	228	10423	0.100	ng	93
31) Chrysene	21.47	228	10749	0.097	ng	92
32) Bis(2-ethylhexyl)phthalate	21.35	149	16416	0.324	ng	# 66
33) Indeno(1,2,3-cd)pyrene	26.61	276	10381	0.025	ng	# 92
35) Benzo(b)fluoranthene	23.18	252	10683m	0.093	ng	
36) Benzo(k)fluoranthene	23.23	252	10478	0.095	ng	97
37) Benzo(a)pyrene	23.84	252	9367	0.095	ng	# 94
38) Dibenzo(a,h)anthracene	26.64	278	9248	0.090	ng	# 90
39) Benzo(g,h,i)perylene	27.43	276	9707	0.023	ng	# 91

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

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