

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE051518\
 Data File : BE096577.D
 Acq On : 15 May 2018 13:45
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE051518

Quant Time: Jun 01 05:04:52 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 15 13:02:03 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	104	0.00
2	1,4-Dioxane	0.600	0.590	1.7	115	0.00
3	n-Nitrosodimethylamine	0.750	0.727	3.1	110	0.00
4 S	2-Fluorophenol	1.103	1.074	2.6	113	-0.01
5 S	Phenol-d6	1.561	1.475	5.5	104	-0.01
6	bis(2-Chloroethyl)ether	1.460	1.390	4.8	106	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	107	-0.01
8 S	Nitrobenzene-d5	0.426	0.386	9.4	104	0.00
9	Naphthalene	4.208	3.890	7.6	104	0.00
10	Hexachlorobutadiene	0.235	0.236	-0.4	123	0.00
11 SURR	2-Methylnaphthalene-d10	0.653	0.593	9.2	103	0.00
12	2-Methylnaphthalene	0.721	0.653	9.4	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
14 S	2,4,6-Tribromophenol	0.158	0.150	5.1	101	0.00
15 S	2-Fluorobiphenyl	1.717	1.688	1.7	104	0.00
16	Acenaphthylene	2.088	1.947	6.8	99	-0.01
17	Acenaphthene	1.272	1.213	4.6	99	-0.01
18	Fluorene	1.583	1.824	-15.2	121	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.243	0.221	9.1	97	0.00
21	Hexachlorobenzene	0.253	0.232	8.3	97	0.00
22	Pentachlorophenol	0.053	0.040	24.5	117	0.00
23	Phenanthrene	1.123	1.060	5.6	99	0.00
24	Anthracene	1.045	0.957	8.4	97	0.00
25 SURR	Fluoranthene-d10	4.890	4.532	7.3	99	0.00
26	Fluoranthene	1.343	1.248	7.1	99	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28	Pyrene	0.722	0.650	10.0	94	0.00
29 S	Terphenyl-d14	0.905	0.870	3.9	98	0.00
30	Benzo(a)anthracene	1.230	1.148	6.7	98	0.00
31	Chrysene	1.280	1.244	2.8	101	0.00
32	Bis(2-ethylhexyl)phthalate	2.632	2.343	11.0	97	0.00
33	Indeno(1,2,3-cd)pyrene	1.227	1.168	4.8	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.234	1.160	6.0	100	0.00
36	Benzo(k)fluoranthene	1.306	1.217	6.8	99	0.00
37 C	Benzo(a)pyrene	1.180	1.117	5.3	100	0.00
38	Dibenzo(a,h)anthracene	1.079	1.006	6.8	98	0.00
39	Benzo(g,h,i)perylene	1.129	1.061	6.0	100	0.00

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(#) = Out of Range	SPCC's out = 0 CCC's out = 0			