

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090115\
 Data File : BE090677.D
 Acq On : 1 Sep 2015 21:45
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE090115

Manual Integrations
 APPROVED

MMDadoda
 9/2/2015 4:07:52 PM

Quant Time: Sep 02 11:46:04 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	41850	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	179348	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	87107	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	222494	5.00	ng	0.00
25) Chrysene-d12	21.07	240	292449	5.00	ng	0.00
32) Perylene-d12	23.36	264	264577	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	6881	0.86	ng	0.00
5) Phenol-d6	6.76	99	9230	0.79	ng	0.01
7) Nitrobenzene-d5	8.71	82	9654	0.81	ng	0.00
13) 2,4,6-Tribromophenol	15.69	330	1605	0.77	ng	0.02
14) 2-Fluorobiphenyl	12.80	172	21127	0.88	ng	0.00
27) Terphenyl-d14	19.54	244	25319	0.78	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.16	88	5779	0.92	ng	97
3) n-Nitrosodimethylamine	3.45	42	6783	0.85	ng	# 99
8) Nitrobenzene	8.76	77	10593	0.82	ng	98
9) Naphthalene	10.36	128	34446	0.88	ng	100
10) Hexachlorobutadiene	10.65	225	5630	0.92	ng	100
11) 2-Methylnaphthalene	11.99	142	17975	0.85	ng	100
15) Acenaphthylene	13.89	152	117452	0.84	ng	100
16) Acenaphthene	14.23	154	20947	0.90	ng	99
17) Fluorene	15.22	166	25498	0.88	ng	99
19) 4-Bromophenyl-phenylether	16.13	248	6608	0.87	ng	95
20) Hexachlorobenzene	16.23	284	37571	0.94	ng	97
21) Pentachlorophenol	16.61	266	1449m	0.90	ng	
22) Phenanthrene	16.96	178	48436	0.88	ng	99
23) Anthracene	17.05	178	38476	0.84	ng	100
24) Fluoranthene	18.97	202	46085	0.79	ng	99
26) Pyrene	19.33	202	46369	0.77	ng	99
28) Benzo(a)anthracene	21.06	228	54345	0.83	ng	99
29) Chrysene	21.11	228	65455	0.91	ng	99
30) Bis(2-ethylhexyl)phthalate	21.01	149	11428	0.72	ng	98
31) Indeno(1,2,3-cd)pyrene	25.73	276	54415	0.79	ng	100
33) Benzo(b)fluoranthene	22.67	252	46218	0.81	ng	99
34) Benzo(k)fluoranthene	22.71	252	55943	0.84	ng	99
35) Benzo(a)pyrene	23.26	252	40139	0.78	ng	99
36) Dibenzo(a,h)anthracene	25.74	278	42946	0.81	ng	99
37) Benzo(g,h,i)perylene	26.45	276	47890	0.81	ng	97

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

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