

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072215\
 Data File : BE090225.D
 Acq On : 22 Jul 2015 8:45
 Operator : TP/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4EC

Manual Integrations
 APPROVED

apatel
 7/22/2015 3:27:05 PM

Quant Time: Jul 22 09:18:49 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE072215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 05:40:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	760	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	3672	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	2076	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5401	0.40	ng/ul	0.00
16) Chrysene-d12	21.11	240	3375	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	1771	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	2323	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	6039	0.39	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.41	128	3884	0.40	ng/ul	100
5) 2-Methylnaphthalene	12.03	142	2559	0.39	ng/ul	100
7) Acenaphthylene	13.93	152	17469	0.40	ng/ul	98
8) Acenaphthene	14.27	153	12490	0.39	ng/ul	98
9) Fluorene	15.26	166	3629	0.38	ng/ul	98
11) Pentachlorophenol	16.61	266	106	0.32	ng/ul	92
12) Phenanthrene	16.99	178	23907	0.38	ng/ul	94
13) Anthracene	17.07	178	24125	0.40	ng/ul	99
15) Fluoranthene	19.00	202	27118	0.39	ng/ul	98
17) Pyrene	19.36	202	27626	0.50	ng/ul	100
18) Benzo(a)anthracene	21.10	228	1533	0.29	ng/ul	97
19) Chrysene	21.15	228	6440m	0.47	ng/ul	
21) Benzo(b)fluoranthene	22.72	252	793	0.35	ng/ul	93
22) Benzo(k)fluoranthene	22.76	252	2221m	0.33	ng/ul	
23) Benzo(a)pyrene	23.32	252	1173	0.34	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.81	276	3076	0.30	ng/ul	98
25) Dibenzo(a,h)anthracene	25.83	278	522	0.31	ng/ul	88
26) Benzo(g,h,i)perylene	26.55	276	4422	0.30	ng/ul#	89

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

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