

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072215\
 Data File : BE090217.D
 Acq On : 22 Jul 2015 1:59
 Operator : TP/IZ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Manual Integrations
 APPROVED

apatel
 7/22/2015 3:26:22 PM

Quant Time: Jul 22 05:37:01 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 04:13:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	792	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	3819	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	2179	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5771	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	3997	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	2108	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2391	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	6480	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.41	128	3949	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.03	142	2668	0.39	ng/ul	99
7) Acenaphthylene	13.93	152	18246	0.40	ng/ul	98
8) Acenaphthene	14.27	153	12928	0.39	ng/ul	98
9) Fluorene	15.26	166	3745	0.37	ng/ul	98
11) Pentachlorophenol	16.61	266	106	0.30	ng/ul	94
12) Phenanthrene	16.99	178	25422	0.38	ng/ul	95
13) Anthracene	17.07	178	25288	0.40	ng/ul	99
15) Fluoranthene	19.00	202	28820	0.38	ng/ul	100
17) Pyrene	19.36	202	29354	0.45	ng/ul	100
18) Benzo(a)anthracene	21.10	228	1731	0.28	ng/ul	96
19) Chrysene	21.15	228	7263m	0.45	ng/ul	
21) Benzo(b)fluoranthene	22.72	252	1000	0.37	ng/ul	96
22) Benzo(k)fluoranthene	22.77	252	2637m	0.33	ng/ul	
23) Benzo(a)pyrene	23.32	252	1456	0.35	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.82	276	4030	0.33	ng/ul	97
25) Dibenzo(a,h)anthracene	25.84	278	620	0.31	ng/ul	92
26) Benzo(g,h,i)perylene	26.56	276	5433	0.31	ng/ul	91

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

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