

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088473.D
 Acq On : 2 Nov 2014 3:40
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Manual Integrations
 APPROVED

mohammad
 11/4/2014 9:08:14 AM

Quant Time: Nov 02 05:44:38 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE103114.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 01 00:42:07 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.09	152	25018	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	96304	5.00	ng	0.00
7) Acenaphthene-d10	10.80	164	44366	5.00	ng	0.01
12) Phenanthrene-d10	12.91	188	85520	5.00	ng	0.00
16) Chrysene-d12	18.33	240	52493	5.00	ng	0.02
22) Perylene-d12	21.40	264	52452	5.00	ng	0.02
System Monitoring Compounds						
4) Nitrobenzene-d5	7.77	82	7435	1.05	ng	0.00
8) 2-Fluorobiphenyl	9.99	172	12973	1.04	ng	0.00
18) Terphenyl-d14	16.08	244	9409	0.98	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.30	88	3289	1.00	ng	100
5) Naphthalene	8.68	128	20835	1.05	ng	99
6) 2-Methylnaphthalene	9.53	142	12860	1.04	ng	97
9) Acenaphthylene	10.62	152	74915	1.04	ng	100
10) Acenaphthene	10.84	154	10713	1.01	ng	99
11) Fluorene	11.49	166	12082	1.07	ng	100
13) Phenanthrene	12.95	178	83161	1.07	ng	97
14) Anthracene	13.03	178	65812	1.05	ng	98
15) Fluoranthene	15.19	202	14688	1.16	ng	100
17) Pyrene	15.65	202	15423	0.90	ng	100
19) Benzo(a)anthracene	18.31	228	53981	1.09	ng	99
20) Chrysene	18.38	228	50351	1.04	ng	100
21) Indeno(1,2,3-cd)pyrene	23.58	276	60836m	1.35	ng	
23) Benzo(b)fluoranthene	20.64	252	12904	1.12	ng	97
24) Benzo(k)fluoranthene	20.69	252	13022	1.03	ng	96
25) Benzo(a)pyrene	21.28	252	12548	1.14	ng	96
26) Dibenzo(a,h)anthracene	23.68	278	11940	1.16	ng	96
27) Benzo(g,h,i)perylene	24.21	276	52225	1.11	ng	99

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

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