

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102114\
 Data File : BE088166.D
 Acq On : 22 Oct 2014 5:47
 Operator : TP/JJ
 Sample : F4303-14DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YS19-SB04-1014DL

Quant Time: Oct 22 06:47:43 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 21 02:11:14 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	22265	5.00	ng	0.00
3) Naphthalene-d8	8.70	136	87594	5.00	ng	0.00
7) Acenaphthene-d10	10.85	164	47456	5.00	ng	0.00
12) Phenanthrene-d10	12.96	188	90888	5.00	ng	0.00
16) Chrysene-d12	18.38	240	52385	5.00	ng	0.00
22) Perylene-d12	21.44	264	43814	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.83	82	82478	13.06	ng	0.00
8) 2-Fluorobiphenyl	10.04	172	132465	10.11	ng	0.00
18) Terphenyl-d14	16.15	244	103789	12.31	ng	0.00
Target Compounds						Qvalue
9) Acenaphthylene	10.67	152	5219	0.07	ng	99
10) Acenaphthene	10.88	154	1563	0.14	ng	88
11) Fluorene	11.55	166	1718	0.15	ng	97
13) Phenanthrene	13.01	178	136434	1.90	ng	95
14) Anthracene	13.09	178	29190	0.41	ng	97
15) Fluoranthene	15.25	202	56476	3.27	ng	99
17) Pyrene	15.70	202	49407	3.47	ng	96
19) Benzo(a)anthracene	18.36	228	80366	1.73	ng	98
20) Chrysene	18.44	228	84081	1.84	ng	97
21) Indeno(1,2,3-cd)pyrene	23.63	276	32385	0.64	ng	# 89
23) Benzo(b)fluoranthene	20.69	252	17713	1.70	ng	98
24) Benzo(k)fluoranthene	20.74	252	7544	0.68	ng	97
25) Benzo(a)pyrene	21.32	252	12673	1.32	ng	98
26) Dibenzo(a,h)anthracene	23.72	278	1781	0.19	ng	# 84
27) Benzo(g,h,i)perylene	24.26	276	30885	0.74	ng	98

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

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