

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090415\
 Data File : BE090732.D
 Acq On : 4 Sep 2015 16:38
 Operator : UM/IZ
 Sample : PB85414BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB85414BS

Quant Time: Sep 04 17:26:12 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	7.55	152	41372	5.00	ng	0.00
6) Naphthalene-d8	10.30	136	188305	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	102062	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	246413	5.00	ng	0.00
25) Chrysene-d12	21.07	240	324625	5.00	ng	0.00
32) Perylene-d12	23.36	264	293766	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.18	112	76420	9.47	ng	0.00
5) Phenol-d6	6.74	99	112399	9.76	ng	0.00
7) Nitrobenzene-d5	8.69	82	54670	4.39	ng	-0.01
13) 2,4,6-Tribromophenol	15.67	330	31300	10.40	ng	0.00
14) 2-Fluorobiphenyl	12.79	172	112125	3.97	ng	-0.01
27) Terphenyl-d14	19.53	244	173187	4.82	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.15	88	22456	4.15	ng	98
3) n-Nitrosodimethylamine	3.45	42	26933	3.66	ng	# 88
8) Nitrobenzene	8.73	77	57348	3.78	ng	98
9) Naphthalene	10.35	128	147865	3.60	ng	99
10) Hexachlorobutadiene	10.65	225	21928	3.41	ng	99
11) 2-Methylnaphthalene	11.97	142	96910	4.36	ng	99
15) Acenaphthylene	13.88	152	663265	4.07	ng	100
16) Acenaphthene	14.23	154	102081	3.75	ng	90
17) Fluorene	15.22	166	136125	4.00	ng	96
19) 4-Bromophenyl-phenylether	16.11	248	34254	4.08	ng	98
20) Hexachlorobenzene	16.23	284	167993	4.17	ng	97
21) Pentachlorophenol	16.58	266	34544	9.98	ng	96
22) Phenanthrene	16.94	178	235097	3.85	ng	99
23) Anthracene	17.05	178	229248	4.50	ng	100
24) Fluoranthene	18.96	202	278886	4.34	ng	99
26) Pyrene	19.32	202	308412	4.59	ng	99
28) Benzo(a)anthracene	21.05	228	301297	4.23	ng	98
29) Chrysene	21.11	228	308946	4.11	ng	99
30) Bis(2-ethylhexyl)phthalate	21.00	149	123398	5.02	ng	99
31) Indeno(1,2,3-cd)pyrene	25.72	276	329916	4.30	ng	99
33) Benzo(b)fluoranthene	22.66	252	275143	4.32	ng	99
34) Benzo(k)fluoranthene	22.71	252	315463	4.25	ng	100
35) Benzo(a)pyrene	23.25	252	268024	4.70	ng	99
36) Dibenzo(a,h)anthracene	25.74	278	271552	3.94	ng	99
37) Benzo(g,h,i)perylene	26.44	276	283337	4.32	ng	98

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

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