

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021518\
 Data File : BF103024.D
 Acq On : 15 Feb 2018 18:16
 Operator : SJ/JU
 Sample : J1540-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CB-3-LINE-NW@3

Manual Integrations
 APPROVED

Sohil
 2/16/2018 2:31:17 PM

Quant Time: Feb 16 00:46:31 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 15 14:08:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	178111	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	726724	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	307682	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	408636	20.00	ng	0.01
75) Chrysene-d12	14.09	240	161192	20.00	ng	0.04
86) Perylene-d12	15.59	264	201110	20.00	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1208218	103.02	ng	0.00
7) Phenol-d6	6.51	99	1455054	103.04	ng	0.00
23) Nitrobenzene-d5	7.44	82	920371	87.86	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	232158	93.75	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1358343	73.26	ng	0.00
78) Terphenyl-d14	13.00	244	858055	126.04	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.52	94	130552	8.627	ng	99
31) Naphthalene	8.18	128	520540	14.661	ng	99
37) 2-Methylnaphthalene	8.87	142	227060	9.768	ng	98
45) 1,1'-Biphenyl	9.33	154	94703	3.654	ng	97
48) Acenaphthylene	9.77	152	83654	2.640	ng	# 96
49) Dimethylphthalate	9.63	163	127534	5.316	ng	99
51) Acenaphthene	9.95	154	1010561	53.326	ng	99
54) Dibenzofuran	10.12	168	1008148	37.749	ng	97
57) Fluorene	10.47	166	1271609	65.443	ng	100
70) Phenanthrene	11.46	178	5715648	251.145	ng	99
71) Anthracene	11.50	178	2551502	110.229	ng	99
72) Carbazole	11.64	167	1461075	64.429	ng	98
74) Fluoranthene	12.66	202	5982698	261.281	ng	99
77) Pyrene	12.89	202	4875527	385.723	ng	97
80) Benzo(a)anthracene	14.07	228	2963346	292.316	ng	98
82) Chrysene	14.12	228	2451961m	239.940	ng	
85) Indeno(1,2,3-cd)pyrene	17.13	276	1591519	187.779	ng	94
87) Benzo(b)fluoranthene	15.16	252	3051124m	234.005	ng	
88) Benzo(k)fluoranthene	15.18	252	837570m	67.230	ng	
89) Benzo(a)pyrene	15.54	252	2058367	175.383	ng	95
90) Dibenzo(a,h)anthracene	17.13	278	397805m	41.760	ng	
91) Benzo(g,h,i)perylene	17.59	276	1563147	167.647	ng	97

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

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