

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031517\
 Data File : BF093687.D
 Acq On : 16 Mar 2017 00:05
 Operator : SJ/MA
 Sample : I2264-13 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-DD9-BASE

Quant Time: Mar 16 03:35:28 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 15:05:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	152965	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	538270	20.00	ng	-0.01
38) Acenaphthene-d10	9.75	164	203517	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	320620	20.00	ng	0.00
75) Chrysene-d12	13.89	240	345367	20.00	ng	0.00
86) Perylene-d12	15.29	264	291507	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	664093	72.26	ng	-0.01
7) Phenol-d6	6.38	99	682408	58.88	ng	0.00
23) Nitrobenzene-d5	7.28	82	400974	41.33	ng	-0.02
41) 2,4,6-Tribromophenol	10.56	330	78451	59.17	ng	0.01
44) 2-Fluorobiphenyl	9.08	172	495407	45.28	ng	-0.01
78) Terphenyl-d14	12.82	244	350844	26.41	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.39	94	47104	3.32	ng	94
31) Naphthalene	8.02	128	69230	2.57	ng	98
49) Dimethylphthalate	9.49	163	81462	6.04	ng	99
51) Acenaphthene	9.78	154	40510	3.66	ng	91
54) Dibenzofuran	9.97	168	31308	2.26	ng	92
57) Fluorene	10.31	166	39481m	3.36	ng	
70) Phenanthrene	11.27	178	553234	30.23	ng	97
71) Anthracene	11.33	178	130940	7.07	ng	98
72) Carbazole	11.50	167	58179	3.35	ng	96
74) Fluoranthene	12.46	202	773686	43.77	ng	96
77) Pyrene	12.69	202	795331	31.66	ng	98
80) Benzo(a)anthracene	13.88	228	420959m	20.64	ng	
82) Chrysene	13.91	228	411270m	21.79	ng	
83) Bis(2-ethylhexyl)phthalate	13.85	149	93366	6.34	ng	93
85) Indeno(1,2,3-cd)pyrene	16.65	276	218402	13.83	ng	98
87) Benzo(b)fluoranthene	14.91	252	564394m	31.79	ng	
88) Benzo(k)fluoranthene	14.92	252	134700m	7.87	ng	
89) Benzo(a)pyrene	15.24	252	383559	23.64	ng	96
90) Dibenzo(a,h)anthracene	16.64	278	46082	3.43	ng	# 79
91) Benzo(a,h,i)perylene	17.05	276	205831	14.94	ng	# 85

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

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