

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
 Data File : BF103858.D
 Acq On : 22 Mar 2018 18:05
 Operator : JU/SJ
 Sample : J1993-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6-MH-A

Manual Integrations
 APPROVED

Sohil
 3/23/2018 2:46:46 PM

Quant Time: Mar 23 12:09:37 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 22 11:56:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	159260	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	620134	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	266151	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	384450	20.00	ng	0.00
75) Chrysene-d12	14.16	240	323984	20.00	ng	0.00
86) Perylene-d12	15.70	264	238171	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	949842	90.71	ng	0.00
7) Phenol-d6	6.60	99	1296141	101.18	ng	0.00
23) Nitrobenzene-d5	7.54	82	765570	86.09	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	250141	109.31	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1193745	71.55	ng	0.00
78) Terphenyl-d14	13.10	244	852878	61.10	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.72	163	120576	5.938	ng	Qvalue # 99
70) Phenanthrene	11.53	178	219950	10.459	ng	99
71) Anthracene	11.58	178	43267	2.026	ng	99
74) Fluoranthene	12.74	202	377382	17.418	ng	99
77) Pyrene	12.96	202	359268	15.100	ng	99
80) Benzo(a)anthracene	14.15	228	172901m	8.431	ng	
82) Chrysene	14.19	228	211633m	10.826	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	64614	3.785	ng	95
87) Benzo(b)fluoranthene	15.24	252	200147m	13.301	ng	
88) Benzo(k)fluoranthene	15.27	252	57053m	3.944	ng	
89) Benzo(a)pyrene	15.63	252	127899	9.371	ng	# 94
91) Benzo(g,h,i)perylene	17.76	276	64254	5.589	ng	96

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

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