

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF032218\
 Data File : BF103860.D
 Acq On : 22 Mar 2018 18:59
 Operator : JU/SJ
 Sample : J1991-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WES-0-6

Manual Integrations
 APPROVED

Sohil
 3/23/2018 2:46:49 PM

Quant Time: Mar 23 11:41:07 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	141546	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	551909	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	220089	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	330875	20.00	ng	0.00
75) Chrysene-d12	14.16	240	292928	20.00	ng	0.00
86) Perylene-d12	15.70	264	241960	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	933917	100.35	ng	0.00
7) Phenol-d6	6.60	99	1127066	98.99	ng	0.00
23) Nitrobenzene-d5	7.53	82	672568	84.98	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	204253	107.94	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1092380	79.17	ng	0.00
78) Terphenyl-d14	13.09	244	839658	66.53	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.72	163	87981	5.240	ng	99
70) Phenanthrene	11.53	178	122329	6.759	ng	99
74) Fluoranthene	12.73	202	189785	10.178	ng	95
77) Pyrene	12.96	202	197994	9.204	ng	99
80) Benzo(a)anthracene	14.15	228	90993m	4.907	ng	
82) Chrysene	14.19	228	85722	4.850	ng	98
87) Benzo(b)fluoranthene	15.24	252	77231m	5.052	ng	
89) Benzo(a)pyrene	15.63	252	54651	3.942	ng	# 92
91) Benzo(g,h,i)perylene	17.76	276	27302	2.338	ng	96

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

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