

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103463.D
 Acq On : 6 Mar 2018 18:50
 Operator : SJ/JU
 Sample : J1722-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-15

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:05 PM

Quant Time: Mar 07 00:43:13 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	176582	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	731627	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	311179	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	467292	20.00	ng	0.00
75) Chrysene-d12	14.06	240	292922	20.00	ng	0.00
86) Perylene-d12	15.57	264	261705	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1079425	92.45	ng	0.00
7) Phenol-d6	6.51	99	1294212	90.59	ng	0.00
23) Nitrobenzene-d5	7.44	82	839344	65.52	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	156442	59.39	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1047813	52.80	ng	0.00
78) Terphenyl-d14	12.99	244	690624	56.68	ng	0.00
Target Compounds						
10) Phenol	6.52	94	89368	5.388	ng	Qvalue # 70
49) Dimethylphthalate	9.62	163	86547	3.467	ng	99
70) Phenanthrene	11.43	178	158098	6.148	ng	98
74) Fluoranthene	12.63	202	316040	12.230	ng	98
77) Pyrene	12.86	202	306095	13.403	ng	98
80) Benzo(a)anthracene	14.05	228	138143	7.268	ng	98
82) Chrysene	14.09	228	123458	6.690	ng	98
85) Indeno(1,2,3-cd)pyrene	17.12	276	50112	3.430	ng	97
87) Benzo(b)fluoranthene	15.13	252	139272m	8.094	ng	
88) Benzo(k)fluoranthene	15.16	252	54799m	3.382	ng	
89) Benzo(a)pyrene	15.51	252	98634	6.419	ng	# 93
91) Benzo(a,h,i)perylene	17.59	276	49105	4.232	ng	96

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

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