

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103465.D
 Acq On : 6 Mar 2018 19:43
 Operator : SJ/JU
 Sample : J1722-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-21

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:19 PM

Quant Time: Mar 07 01:00:05 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	155570	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	689010	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	288212	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	436691	20.00	ng	0.00
75) Chrysene-d12	14.06	240	289227	20.00	ng	0.00
86) Perylene-d12	15.58	264	227575	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1025001	99.65	ng	0.00
7) Phenol-d6	6.51	99	1204919	95.73	ng	0.00
23) Nitrobenzene-d5	7.44	82	796883	66.05	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	143071	58.65	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1007026	55.20	ng	0.00
78) Terphenyl-d14	12.99	244	701238	58.28	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	70835	3.064	ng	Qvalue 100
70) Phenanthrene	11.43	178	204848	8.524	ng	99
71) Anthracene	11.49	178	69735	2.830	ng	99
74) Fluoranthene	12.63	202	426516	17.661	ng	99
77) Pyrene	12.86	202	441544	19.581	ng	99
80) Benzo(a)anthracene	14.05	228	190106	10.130	ng	98
82) Chrysene	14.09	228	183507	10.071	ng	97
85) Indeno(1,2,3-cd)pyrene	17.13	276	62285	4.318	ng	97
87) Benzo(b)fluoranthene	15.13	252	171300m	11.448	ng	
88) Benzo(k)fluoranthene	15.16	252	62814m	4.458	ng	
89) Benzo(a)pyrene	15.52	252	118729	8.886	ng	# 93
91) Benzo(a,h,i)perylene	17.60	276	63469	6.290	ng	93

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

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