

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103717.D
 Acq On : 17 Mar 2018 00:40
 Operator : JU/SJ
 Sample : J1945-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-A

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:02:47 PM

Quant Time: Mar 17 04:06:54 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	172281	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	688620	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	296078	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	437082	20.00	ng	0.00
75) Chrysene-d12	14.15	240	324342	20.00	ng	0.00
86) Perylene-d12	15.69	264	256991	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1159557	102.37	ng	0.01
7) Phenol-d6	6.59	99	1407977	101.60	ng	0.00
23) Nitrobenzene-d5	7.53	82	842530	85.32	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	256232	100.65	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1237400	66.67	ng	0.00
78) Terphenyl-d14	13.09	244	893922	63.97	ng	0.00
Target Compounds						
10) Phenol	6.60	94	37692	2.560	ng	Qvalue # 56
49) Dimethylphthalate	9.72	163	144492	6.397	ng	100
70) Phenanthrene	11.53	178	162543	6.798	ng	97
74) Fluoranthene	12.72	202	200129	8.125	ng	99
77) Pyrene	12.95	202	156608	6.575	ng	99
80) Benzo(a)anthracene	14.14	228	64146	3.124	ng	99
82) Chrysene	14.17	228	66063	3.376	ng	98
87) Benzo(b)fluoranthene	15.23	252	70963m	4.371	ng	
89) Benzo(a)pyrene	15.62	252	42134	2.861	ng	# 92

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

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