

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090717\
 Data File : BF098335.D
 Acq On : 8 Sep 2017 1:18
 Operator : SJ/JU
 Sample : I5071-14
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G55A-2.5-4.5

Manual Integrations
 APPROVED

Sohil
 9/8/2017 5:45:31 PM

Quant Time: Sep 08 04:13:33 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	109978	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	413775	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	171192	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	275458	20.00	ng	-0.01
75) Chrysene-d12	13.84	240	234933	20.00	ng	-0.02
86) Perylene-d12	15.26	264	193659	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	819888	113.40	ng	0.00
7) Phenol-d6	6.34	99	1135200	115.56	ng	-0.01
23) Nitrobenzene-d5	7.26	82	665310	87.35	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	175398	118.08	ng	-0.02
44) 2-Fluorobiphenyl	9.05	172	1007583	86.47	ng	-0.02
78) Terphenyl-d14	12.80	244	686202	60.91	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	171476	13.703	ng	99
70) Phenanthrene	11.23	178	63952	4.357	ng	99
74) Fluoranthene	12.42	202	123507	8.061	ng	99
77) Pyrene	12.65	202	122658	6.384	ng	98
80) Benzo(a)anthracene	13.83	228	86264	6.126	ng	93
82) Chrysene	13.87	228	91600	6.540	ng	96
83) Bis(2-ethylhexyl)phthalate	13.83	149	27791	3.404	ng	# 99
85) Indeno(1,2,3-cd)pyrene	16.62	276	33639	3.265	ng	93
87) Benzo(b)fluoranthene	14.86	252	141542m	11.595	ng	
88) Benzo(k)fluoranthene	14.88	252	35499m	3.095	ng	
89) Benzo(a)pyrene	15.20	252	68675	6.355	ng	99
91) Benzo(a,h,i)perylene	17.02	276	32679	3.560	ng	97

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

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