

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104162.D
 Acq On : 3 Apr 2018 5:43
 Operator : JU/SJ
 Sample : J2121-14
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PA-10-COMP

Manual Integrations
 APPROVED

Sohil
 4/3/2018 2:28:10 PM

Quant Time: Apr 03 12:58:41 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 02 13:40:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	98127	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	395789	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	168320	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	247916	20.00	ng	0.00
75) Chrysene-d12	14.15	240	206712	20.00	ng	0.00
86) Perylene-d12	15.70	264	135007	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	542340	88.14	ng	0.01
7) Phenol-d6	6.60	99	718175	94.87	ng	0.00
23) Nitrobenzene-d5	7.52	82	400714	61.92	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	141455	75.47	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	739923	69.32	ng	0.00
78) Terphenyl-d14	13.08	244	521157	58.21	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	44453	3.345	ng	Qvalue 100
70) Phenanthrene	11.52	178	40926	3.049	ng	99
74) Fluoranthene	12.72	202	84021	5.889	ng	96
77) Pyrene	12.94	202	96871	6.519	ng	99
80) Benzo(a)anthracene	14.14	228	44823	3.465	ng	# 89
82) Chrysene	14.18	228	57421m	4.533	ng	
85) Indeno(1,2,3-cd)pyrene	17.30	276	50003	3.809	ng	# 88
87) Benzo(b)fluoranthene	15.24	252	80645m	9.815	ng	
88) Benzo(k)fluoranthene	15.26	252	33501m	4.328	ng	
89) Benzo(a)pyrene	15.63	252	66863	8.781	ng	# 86
91) Benzo(g,h,i)perylene	17.79	276	44580	6.322	ng	# 90

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

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