

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053017\
 Data File : BF095547.D
 Acq On : 30 May 2017 21:29
 Operator : SJ/MA
 Sample : I3302-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM17-COMP-0-6

Manual Integrations
 APPROVED

mohammad
 5/31/2017 5:02:53 PM

Quant Time: May 31 04:35:51 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 26 16:50:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	94338	20.00	ng	0.00
21) Naphthalene-d8	9.77	136	340930	20.00	ng	0.00
38) Acenaphthene-d10	12.59	164	136557	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	203747	20.00	ng	0.00
75) Chrysene-d12	18.67	240	206642	20.00	ng	0.00
86) Perylene-d12	20.35	264	141727	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	520157	91.93	ng	0.00
7) Phenol-d6	7.30	99	613324	90.34	ng	0.00
23) Nitrobenzene-d5	8.65	82	326822	60.45	ng	0.00
41) 2,4,6-Tribromophenol	13.91	330	86475	77.32	ng	0.00
44) 2-Fluorobiphenyl	11.54	172	642395	67.71	ng	0.00
78) Terphenyl-d14	17.31	244	407767	43.46	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.24	163	93573	8.35	ng	98
70) Phenanthrene	15.04	178	66024	5.64	ng	96
74) Fluoranthene	16.78	202	148820	12.39	ng	98
77) Pyrene	17.08	202	138217	8.94	ng	97
80) Benzo(a)anthracene	18.66	228	74213	6.17	ng	96
82) Chrysene	18.71	228	73165	6.29	ng	96
85) Indeno(1,2,3-cd)pyrene	21.70	276	28591	3.03	ng	93
87) Benzo(b)fluoranthene	19.95	252	80485m	9.19	ng	
88) Benzo(k)fluoranthene	19.97	252	28119m	3.33	ng	
89) Benzo(a)pyrene	20.30	252	52111	6.45	ng	94
91) Benzo(g,h,i)perylene	22.09	276	30089	4.59	ng	92

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

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