

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053017\
 Data File : BF095546.D
 Acq On : 30 May 2017 20:57
 Operator : SJ/MA
 Sample : I3302-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM16-COMP-6-18

Manual Integrations
 APPROVED

mohammad
 5/31/2017 5:02:51 PM

Quant Time: May 31 04:31:24 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	94595	20.00	ng	0.00
21) Naphthalene-d8	9.77	136	362544	20.00	ng	0.00
38) Acenaphthene-d10	12.60	164	143030	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	210903	20.00	ng	0.00
75) Chrysene-d12	18.67	240	214443	20.00	ng	0.00
86) Perylene-d12	20.35	264	150417	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	629236	110.90	ng	0.00
7) Phenol-d6	7.29	99	714137	104.90	ng	0.00
23) Nitrobenzene-d5	8.64	82	372937	64.87	ng	-0.01
41) 2,4,6-Tribromophenol	13.90	330	106935	91.29	ng	0.00
44) 2-Fluorobiphenyl	11.53	172	686224	69.06	ng	0.00
78) Terphenyl-d14	17.31	244	405722	41.67	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.24	163	96973	8.26	ng	100
70) Phenanthrene	15.04	178	38266	3.16	ng	96
74) Fluoranthene	16.78	202	85248	6.86	ng	99
77) Pyrene	17.09	202	81368	5.07	ng	97
80) Benzo(a)anthracene	18.66	228	39167m	3.14	ng	
82) Chrysene	18.71	228	45144	3.74	ng	98
87) Benzo(b)fluoranthene	19.95	252	42089m	4.53	ng	
89) Benzo(a)pyrene	20.30	252	30162	3.52	ng	94
91) Benzo(g,h,i)perylene	22.09	276	16266	2.34	ng	96

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

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