

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053017\
 Data File : BF095533.D
 Acq On : 30 May 2017 13:59
 Operator : SJ/MA
 Sample : I3281-13
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R1-TP7-D34

Manual Integrations
 APPROVED

mohammad
 5/31/2017 5:02:18 PM

Quant Time: May 31 02:18:14 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	94117	20.00	ng	0.00
21) Naphthalene-d8	9.77	136	391916	20.00	ng	0.00
38) Acenaphthene-d10	12.60	164	190612	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	333390	20.00	ng	0.00
75) Chrysene-d12	18.67	240	193355	20.00	ng	0.00
86) Perylene-d12	20.35	264	163744	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	620884	109.98	ng	0.00
7) Phenol-d6	7.28	99	768322	113.43	ng	-0.01
23) Nitrobenzene-d5	8.64	82	452846	72.86	ng	-0.01
41) 2,4,6-Tribromophenol	13.89	330	160405	102.75	ng	0.00
44) 2-Fluorobiphenyl	11.53	172	829835	62.66	ng	0.00
78) Terphenyl-d14	17.31	244	581252	66.21	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.24	163	202734	12.96	ng	99
70) Phenanthrene	15.04	178	70851	3.70	ng	98
74) Fluoranthene	16.78	202	142394	7.25	ng	97
77) Pyrene	17.08	202	117304	8.11	ng	99
80) Benzo(a)anthracene	18.66	228	57150	5.08	ng	95
82) Chrysene	18.71	228	67063m	6.16	ng	
85) Indeno(1,2,3-cd)pyrene	21.70	276	25188	2.85	ng	94
87) Benzo(b)fluoranthene	19.95	252	57464m	5.68	ng	
88) Benzo(k)fluoranthene	19.98	252	22854m	2.34	ng	
89) Benzo(a)pyrene	20.30	252	41594	4.46	ng	94
91) Benzo(g,h,i)perylene	22.09	276	24771	3.27	ng	98

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

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