

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060617\
 Data File : BF095727.D
 Acq On : 7 Jun 2017 00:48
 Operator : SJ/MA
 Sample : I3482-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-060517-A

Manual Integrations
 APPROVED

mohammad
 6/7/2017 2:18:57 PM

Quant Time: Jun 07 04:22:09 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	80264	20.00	ng	-0.02
21) Naphthalene-d8	9.41	136	299731	20.00	ng	-0.02
38) Acenaphthene-d10	12.23	164	120033	20.00	ng	-0.02
63) Phenanthrene-d10	14.63	188	182207	20.00	ng	-0.02
75) Chrysene-d12	18.34	240	153414	20.00	ng	-0.01
86) Perylene-d12	20.01	264	107710	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	516053	102.45	ng	-0.01
7) Phenol-d6	6.94	99	590822	97.98	ng	-0.02
23) Nitrobenzene-d5	8.30	82	334551	69.32	ng	-0.02
41) 2,4,6-Tribromophenol	13.53	330	90475	85.19	ng	-0.02
44) 2-Fluorobiphenyl	11.19	172	579249	67.15	ng	-0.02
78) Terphenyl-d14	17.00	244	360562	58.33	ng	-0.02
Target Compounds						
49) Dimethylphthalate	11.89	163	93475	10.257	ng	Qvalue # 98
70) Phenanthrene	14.66	178	68392	6.505	ng	98
74) Fluoranthene	16.44	202	108820	10.458	ng	98
77) Pyrene	16.74	202	96640	8.442	ng	# 92
80) Benzo(a)anthracene	18.33	228	43328	4.858	ng	94
82) Chrysene	18.37	228	39208m	4.399	ng	
87) Benzo(b)fluoranthene	19.61	252	38020m	5.637	ng	
88) Benzo(k)fluoranthene	19.64	252	14895m	2.311	ng	
89) Benzo(a)pyrene	19.96	252	27127	4.376	ng	# 96
91) Benzo(g,h,i)perylene	21.53	276	15252	2.845	ng	# 76

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

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