

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085604.D
 Acq On : 20 Mar 2016 8:02
 Operator : UM/SJ
 Sample : H1815-16
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMP

Manual Integrations
 APPROVED

apatel
 3/21/2016 5:10:45 PM

Quant Time: Mar 21 04:58:50 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	135346	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	579702	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	247996	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	478954	20.00	ng	-0.01
75) Chrysene-d12	14.00	240	295459	20.00	ng	-0.01
86) Perylene-d12	15.43	264	261339	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1109307	137.67	ng	0.00
7) Phenol-d6	6.49	99	1501130	145.25	ng	0.00
23) Nitrobenzene-d5	7.42	82	876021	87.72	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	367166	151.14	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1334876	94.45	ng	-0.01
78) Terphenyl-d14	12.96	244	1194485	97.28	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	152568	8.15	ng	99
70) Phenanthrene	11.40	178	70953	2.81	ng	98
74) Fluoranthene	12.59	202	130469	4.93	ng	90
77) Pyrene	12.81	202	131138	6.18	ng	99
80) Benzo(a)anthracene	13.99	228	58919m	3.44	ng	
82) Chrysene	14.04	228	49990	3.03	ng	97
85) Indeno(1,2,3-cd)pyrene	16.83	276	34748	2.52	ng	94
87) Benzo(b)fluoranthene	15.02	252	74337m	4.36	ng	
89) Benzo(a)pyrene	15.37	252	52091	3.60	ng	96
91) Benzo(g,h,i)perylene	17.25	276	35734	3.06	ng	96

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

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