

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
 Data File : BF078948.D
 Acq On : 5 May 2015 18:54
 Operator : TP/IZ
 Sample : G2044-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-49S

Quant Time: May 06 02:07:41 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 00:45:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.36	152	77695	20.00	ng	0.00
21) Naphthalene-d8	8.94	136	310853	20.00	ng	0.00
38) Acenaphthene-d10	11.11	164	151644	20.00	ng	0.00
63) Phenanthrene-d10	12.95	188	284164	20.00	ng	-0.01
75) Chrysene-d12	16.24	240	308663	20.00	ng	-0.03
86) Perylene-d12	17.95	264	336152	20.00	ng	-0.14

System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	415837	92.13	ng	0.01
7) Phenol-d6	6.90	99	528753	88.63	ng	0.00
23) Nitrobenzene-d5	8.04	82	321223	59.39	ng	0.00
41) 2,4,6-Tribromophenol	12.09	330	140369	85.56	ng	0.00
44) 2-Fluorobiphenyl	10.27	172	576382	52.95	ng	-0.01
78) Terphenyl-d14	14.95	244	620661	48.14	ng	0.00

Target Compounds						Qvalue
49) Dimethylphthalate	10.78	163	22265	2.00	ng	99
70) Phenanthrene	12.98	178	236187	14.86	ng	99
71) Anthracene	13.04	178	60039	3.85	ng	99
74) Fluoranthene	14.46	202	391428	23.63	ng	99
77) Pyrene	14.74	202	367722	20.67	ng	98
80) Benzo(a)anthracene	16.23	228	225813	13.36	ng	97
82) Chrysene	16.27	228	179569	11.05	ng	96
83) Bis(2-ethylhexyl)phthalate	16.29	149	26333	2.24	ng	95
85) Indeno(1,2,3-cd)pyrene	19.44	276	145731m	7.04	ng	
87) Benzo(b)fluoranthene	17.51	252	265836	14.45	ng	98
88) Benzo(k)fluoranthene	17.53	252	107817m	6.05	ng	
89) Benzo(a)pyrene	17.89	252	206057	12.16	ng	97
91) Benzo(a,h,i)perylene	19.89	276	159327	8.83	ng	99

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

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