

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011515\
 Data File : BF076890.D
 Acq On : 15 Jan 2015 18:33
 Operator : IZ / TP
 Sample : G1030-04
 Misc : S
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2-SS-1(2-4)

Quant Time: Jan 16 01:10:14 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 15 11:52:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	33757	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	144715	20.00	ng	0.00
38) Acenaphthene-d10	15.07	164	72264	20.00	ng	-0.01
63) Phenanthrene-d10	17.82	188	118315	20.00	ng	0.00
75) Chrysene-d12	21.31	240	115009	20.00	ng	0.00
86) Perylene-d12	22.99	264	108306	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.09	112	248291	120.93	ng	0.00
7) Phenol-d6	7.42	99	313157	120.91	ng	0.00
23) Nitrobenzene-d5	9.32	82	190354	72.87	ng	0.00
41) 2,4,6-Tribromophenol	16.72	330	65183	111.12	ng	0.00
44) 2-Fluorobiphenyl	13.58	172	314448	66.22	ng	-0.01
78) Terphenyl-d14	20.04	244	274308	54.91	ng	0.00
Target Compounds						Qvalue
37) 2-Methylnaphthalene	12.63	142	20260	3.98	ng	97
49) Dimethylphthalate	14.64	163	69827	13.63	ng	97
70) Phenanthrene	17.85	178	26208	3.55	ng	96
74) Fluoranthene	19.49	202	56759	7.51	ng	98
77) Pyrene	19.77	202	53046	6.73	ng	# 93
80) Benzo(a)anthracene	21.29	228	32714	4.53	ng	98
82) Chrysene	21.34	228	26851	4.08	ng	98
85) Indeno(1,2,3-cd)pyrene	24.13	276	20252	2.90	ng	# 79
87) Benzo(b)fluoranthene	22.56	252	35996m	4.93	ng	
88) Benzo(k)fluoranthene	22.59	252	13287m	2.08	ng	
89) Benzo(a)pyrene	22.92	252	29393	4.57	ng	94
91) Benzo(a,h,i)perylene	24.43	276	26997	4.60	ng	# 92

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

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