

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\
 Data File : BF076833.D
 Acq On : 13 Jan 2015 7:02
 Operator : IZ / TP
 Sample : G1030-02
 Misc : S
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-SS-2(4-6)

Quant Time: Jan 13 11:38:51 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 10 06:22:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	32103	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	138773	20.00	ng	0.00
38) Acenaphthene-d10	15.13	164	73531	20.00	ng	0.00
63) Phenanthrene-d10	17.87	188	115238	20.00	ng	0.01
75) Chrysene-d12	21.36	240	110419	20.00	ng	0.01
86) Perylene-d12	23.05	264	110838	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	181006	92.74	ng	-0.01
7) Phenol-d6	7.49	99	226695	92.19	ng	-0.01
23) Nitrobenzene-d5	9.37	82	140918	55.25	ng	-0.01
41) 2,4,6-Tribromophenol	16.77	330	48926	77.66	ng	-0.01
44) 2-Fluorobiphenyl	13.65	172	224663	44.46	ng	0.00
78) Terphenyl-d14	20.08	244	184026	35.90	ng	0.01
Target Compounds						Qvalue
10) Phenol	7.51	94	6019	2.26	ng	# 61
31) Naphthalene	11.04	128	125496	17.52	ng	98
37) 2-Methylnaphthalene	12.70	142	34589	6.85	ng	98
48) Acenaphthylene	14.78	152	48245	6.14	ng	98
49) Dimethylphthalate	14.70	163	49835	9.23	ng	99
51) Acenaphthene	15.20	154	109444	24.66	ng	98
54) Dibenzofuran	15.63	168	115296	17.95	ng	97
57) Fluorene	16.32	166	149703	27.61	ng	97
70) Phenanthrene	17.91	178	1643156	227.47	ng	98
71) Anthracene	17.97	178	408641	57.26	ng	99
72) Carbazole	18.25	167	185679	26.95	ng	99
74) Fluoranthene	19.55	202	1727347	231.88	ng	94
77) Pvrene	19.83	202	1483350	183.36	ng	# 94
80) Benzo(a)anthracene	21.35	228	787859	109.93	ng	97
82) Chrysene	21.39	228	640831m	100.59	ng	
85) Indeno(1,2,3-cd)pyrene	24.23	276	319095m	48.17	ng	
87) Benzo(b)fluoranthene	22.63	252	732151m	89.96	ng	
88) Benzo(k)fluoranthene	22.64	252	283283m	44.12	ng	
89) Benzo(a)pvrene	23.00	252	537561	79.71	ng	96
90) Dibenzo(a,h)anthracene	24.24	278	66225m	10.80	ng	
91) Benzo(q,h,i)perylene	24.54	276	278281	44.66	ng	98

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

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