

Data Path : U:\HPCHEM1\BNA F\DATA\BF012518\
 Data File : BF102442.D
 Acq On : 25 Jan 2018 23:45
 Operator : SJ/JU
 Sample : J1236-02RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 33E-6RE

Manual Integrations
 APPROVED

Sohil
 1/26/2018 11:24:45 AM

Quant Time: Jan 26 02:52:06 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 25 13:44:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	149330	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	599133	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	248313	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	356596	20.00	ng	0.00
75) Chrysene-d12	13.88	240	282271	20.00	ng	0.00
86) Perylene-d12	15.31	264	246212	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	215965	21.93	ng	0.01
7) Phenol-d6	6.36	99	949470	79.97	ng	0.00
23) Nitrobenzene-d5	7.29	82	724806	69.52	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	16022	6.51	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1084475	64.22	ng	0.00
78) Terphenyl-d14	12.82	244	745301	59.52	ng	0.00
Target Compounds						
10) Phenol	6.37	94	86831	6.699	ng	95
20) 3+4-Methylphenols	7.14	107	34277	3.250	ng	91
49) Dimethylphthalate	9.47	163	193303	9.418	ng	99
70) Phenanthrene	11.26	178	49226	2.369	ng	99
73) Di-n-butylphthalate	11.80	149	82480	3.189	ng	98
74) Fluoranthene	12.45	202	81048	3.825	ng	97
77) Pyrene	12.68	202	152618	6.993	ng	97
80) Benzo(a)anthracene	13.87	228	46508m	2.676	ng	
82) Chrysene	13.90	228	53299	3.020	ng	97
85) Indeno(1,2,3-cd)pyrene	16.70	276	33434	2.294	ng	96
87) Benzo(b)fluoranthene	14.90	252	65570m	4.109	ng	
89) Benzo(a)pyrene	15.24	252	56664	3.937	ng	# 95
91) Benzo(a,h,i)perylene	17.12	276	44993	3.908	ng	# 87

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

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