

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042221\
 Data File : BF123688.D
 Acq On : 22 Apr 2021 12:51
 Operator : JU/CG
 Sample : M1998-03DL 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3CDL

Manual Integrations
 APPROVED

mohammad
 4/23/2021 8:37:08 AM

Quant Time: Apr 22 13:26:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 20 15:49:40 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	263229	20.00	ng	0.00
21) Naphthalene-d8	8.139	136	1025264	20.00	ng	0.00
39) Acenaphthene-d10	9.898	164	522039	20.00	ng	0.00
64) Phenanthrene-d10	11.386	188	1013594	20.00	ng	0.00
76) Chrysene-d12	14.033	240	667494	20.00	ng	0.00
86) Perylene-d12	15.521	264	614677	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	732009	44.95	ng	0.00
7) Phenol-d6	6.486	99	985090	43.65	ng	0.00
23) Nitrobenzene-d5	7.410	82	689037	30.90	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	186599	31.57	ng	0.00
45) 2-Fluorobiphenyl	9.216	172	1006171	28.73	ng	0.00
79) Terphenyl-d14	12.974	244	1033274	30.09	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.498	94	899138	37.51	ng	97
20) 3+4-Methylphenols	7.257	107	126913	6.62	ng	# 65
75) Fluoranthene	12.615	202	178767	3.07	ng	97
78) Pyrene	12.833	202	168909	3.47	ng	97
84) Bis(2-ethylhexyl)phtha...	14.015	149	79812	2.62	ng	100
88) Benzo(b)fluoranthene	15.080	252	94930m	2.19	ng	
90) Benzo(a)pyrene	15.451	252	70010	2.03	ng	# 94

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

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