

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031921\
 Data File : BG048433.D
 Acq On : 19 Mar 2021 21:07
 Operator : CG/JU
 Sample : M1732-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-3A-(0.5-1.0)

Manual Integrations
 APPROVED

mohammad
 3/22/2021 2:49:48 PM

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Quant Time: Mar 19 22:52:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG031821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 18 16:38:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	48464	20.00	ng	0.00
21) Naphthalene-d8	10.930	136	188893	20.00	ng	# 0.00
39) Acenaphthene-d10	14.749	164	152558	20.00	ng	0.00
64) Phenanthrene-d10	17.499	188	380326	20.00	ng	# 0.00
76) Chrysene-d12	21.794	240	386551	20.00	ng	# 0.00
86) Perylene-d12	25.131	264	396477	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.654	112	275180	93.90	ng	0.00
7) Phenol-d6	7.252	99	391766	91.87	ng	0.00
23) Nitrobenzene-d5	9.273	82	298345	63.15	ng	0.00
42) 2,4,6-Tribromophenol	16.236	330	245475	98.94	ng	0.00
45) 2-Fluorobiphenyl	13.368	172	588019	55.45	ng	0.00
79) Terphenyl-d14	20.107	244	996130	46.47	ng	0.00
Target Compounds						
						Qvalue
50) Dimethylphthalate	14.191	163	29896	2.41	ng	100
75) Fluoranthene	19.549	202	56339	2.11	ng	97
78) Pyrene	19.914	202	95978	3.62	ng	96
81) Benzo(a)anthracene	21.770	228	93481	3.49	ng	98
83) Chrysene	21.841	228	128256	4.99	ng	97
88) Benzo(b)fluoranthene	24.062	252	170483	6.10	ng	# 92
90) Benzo(a)pyrene	24.967	252	94075m	3.69	ng	

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

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