

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051021\
 Data File : BP005479.D
 Acq On : 10 May 2021 21:07
 Operator : CG/JU
 Sample : M2310-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BS-SP2

Manual Integrations
 APPROVED

mohammad
 5/11/2021 2:31:59 PM

Quant Time: May 11 04:23:16 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 07 12:37:06 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	10997	20.00	ng	0.00
21) Naphthalene-d8	10.493	136	29396	20.00	ng	#-0.01
39) Acenaphthene-d10	14.351	164	23004	20.00	ng	0.00
64) Phenanthrene-d10	17.098	188	61034	20.00	ng	#-0.01
76) Chrysene-d12	21.192	240	115046	20.00	ng	#-0.01
86) Perylene-d12	23.504	264	142413	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	55694	97.19	ng	0.00
7) Phenol-d6	6.899	99	57537	80.52	ng	0.00
23) Nitrobenzene-d5	8.869	82	49877	66.78	ng	-0.01
42) 2,4,6-Tribromophenol	15.845	330	54493	83.48	ng	-0.01
45) 2-Fluorobiphenyl	12.969	172	109238	53.57	ng	-0.01
79) Terphenyl-d14	19.669	244	314570	46.44	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	13.822	163	6069	3.06	ng	# 97
71) Phenanthrene	17.139	178	37787	10.90	ng	97
72) Anthracene	17.239	178	12127m	3.52	ng	
75) Fluoranthene	19.116	202	80849	16.95	ng	97
78) Pyrene	19.469	202	77174	11.85	ng	99
81) Benzo(a)anthracene	21.180	228	46168	6.19	ng	97
83) Chrysene	21.227	228	44302m	6.13	ng	
87) Indeno(1,2,3-cd)pyrene	25.898	276	30455	2.54	ng	# 94
88) Benzo(b)fluoranthene	22.804	252	61098	6.44	ng	# 96
89) Benzo(k)fluoranthene	22.845	252	21665m	2.33	ng	
90) Benzo(a)pyrene	23.404	252	45195	5.20	ng	99
92) Benzo(g,h,i)perylene	26.633	276	30472	3.12	ng	# 92

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

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