

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091820\
 Data File : BP003325.D
 Acq On : 18 Sep 2020 17:07
 Operator : CG/JU
 Sample : L4073-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H-1

Manual Integrations
 APPROVED

mohammad
 9/21/2020 10:35:44 AM

Quant Time: Sep 18 18:58:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 18 14:13:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	193942	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	803871	20.00	ng	# 0.00
39) Acenaphthene-d10	14.245	164	504300	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	1076032	20.00	ng	0.00
76) Chrysene-d12	21.104	240	914328	20.00	ng	0.00
86) Perylene-d12	23.315	264	829802	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.216	112	915970	82.47	ng	0.00
7) Phenol-d6	6.799	99	1202254	81.21	ng	0.00
23) Nitrobenzene-d5	8.746	82	786838	57.50	ng	0.00
42) 2,4,6-Tribromophenol	15.745	330	625683	81.34	ng	0.00
45) 2-Fluorobiphenyl	12.863	172	1925263	55.84	ng	0.00
79) Terphenyl-d14	19.586	244	2626858	59.86	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.710	163	429986	10.68	ng	99
71) Phenanthrene	17.045	178	469854	8.03	ng	99
72) Anthracene	17.139	178	146060	2.53	ng	99
75) Fluoranthene	19.027	202	1826244	24.81	ng	99
78) Pyrene	19.380	202	1635816	28.69	ng	100
81) Benzo(a)anthracene	21.092	228	1023403	17.31	ng	99
83) Chrysene	21.139	228	867130	15.27	ng	98
87) Indeno(1,2,3-cd)pyrene	25.562	276	457484	7.27	ng	# 94
88) Benzo(b)fluoranthene	22.657	252	990157	18.90	ng	98
89) Benzo(k)fluoranthene	22.692	252	380929m	7.59	ng	
90) Benzo(a)pyrene	23.221	252	761603	15.41	ng	97
91) Dibenzo(a,h)anthracene	25.562	278	126284m	2.43	ng	
92) Benzo(g,h,i)perylene	26.256	276	412079	7.95	ng	96

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

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