

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091820\
 Data File : BP003329.D
 Acq On : 18 Sep 2020 19:23
 Operator : CG/JU
 Sample : L3871-19
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TR-08-091720

Manual Integrations
 APPROVED

mohammad
 9/21/2020 10:35:52 AM

Quant Time: Sep 19 06:19:23 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	182973	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	745789	20.00	ng	# 0.00
39) Acenaphthene-d10	14.245	164	452424	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	916823	20.00	ng	0.00
76) Chrysene-d12	21.104	240	685812	20.00	ng	0.00
86) Perylene-d12	23.321	264	679059	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.216	112	1018506	97.20	ng	0.00
7) Phenol-d6	6.799	99	1297991	92.93	ng	0.00
23) Nitrobenzene-d5	8.751	82	863176	68.00	ng	0.00
42) 2,4,6-Tribromophenol	15.751	330	543858	78.81	ng	0.00
45) 2-Fluorobiphenyl	12.869	172	1895787	61.28	ng	0.00
79) Terphenyl-d14	19.586	244	2149946	65.32	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.428	128	335741	8.52	ng	100
37) 2-Methylnaphthalene	12.051	142	92641	3.25	ng	98
38) 1-Methylnaphthalene	12.275	142	64568m	2.38	ng	
49) Acenaphthylene	13.969	152	102574	2.40	ng	99
50) Dimethylphthalate	13.710	163	398950	11.04	ng	100
58) Fluorene	15.304	166	98891	3.02	ng	99
71) Phenanthrene	17.045	178	472964	9.48	ng	99
72) Anthracene	17.139	178	166940	3.40	ng	100
75) Fluoranthene	19.027	202	724043	11.54	ng	99
78) Pyrene	19.380	202	743779	17.39	ng	100
81) Benzo(a)anthracene	21.092	228	333531	7.52	ng	95
83) Chrysene	21.139	228	311066	7.30	ng	97
87) Indeno(1,2,3-cd)pyrene	25.562	276	199272	3.87	ng	# 95
88) Benzo(b)fluoranthene	22.657	252	387890m	9.05	ng	
89) Benzo(k)fluoranthene	22.692	252	148790m	3.62	ng	
90) Benzo(a)pyrene	23.221	252	320211	7.92	ng	97
92) Benzo(g,h,i)perylene	26.251	276	188094	4.44	ng	98

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

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